



Certificate of Analysis
Number: 1030-26040367-001A

Houston Laboratories
8820 Interchange Drive
Houston, TX 77054
Phone 713-660-0901

Luke Arnsberger
Dolan Intergration Group
11025 Dover St, Ste 800
Westminister, CO 80021

Sample ID: 26041347-001A
Station Name: DCP Wattenburg Pipeline
Station Number: Spill ID 493177
Station Location: Hwy 52 & CR19 (G) 2/2026
Cylinder No: 043483
Instrument: PIANO, WT
Analyzed: 04/17/2026 01:22:29 by HKP

Report Date: 04/22/2026
Sampled By:
Sample Of: Liquid Spot
Sample Date: 04/01/2026 13:17
Sample Conditions:
Received Date: 04/10/2026
Login Date: 04/13/2026
Method: Proprietary

Analytical Data

Carbon Range	C4-C18, C22	Major Range	C5-C9
Paraffins	29.587 wt%	n-Hexane	12.679 wt%
Isoparaffins	33.559 wt%	Benzene	0.937 wt%
Olefins	0.268 wt%	2,2,4-Trimethylpentane	0.019 wt%
Naphthenics	30.203 wt%	Toluene	3.136 wt%
Aromatics	6.383 wt%	Ethylbenzene	0.212 wt%
Unknowns	ND wt%	m-Xylene	0.935 wt%
		p-Xylene	0.334 wt%
		o-Xylene	0.336 wt%
C1-C29 Total	100.000 wt%	Xylenes	1.605 wt%
Naphthalene	0.002 wt%	n-Heptadecane	0.001 wt%
2-Methylnaphthalene	ND wt%	Pristane	ND wt%
1-Methylnaphthalene	ND wt%	n-Octadecane	0.001 wt%
Oxygenates	ND wt%	Phytane	ND wt%
		Ethanol	ND wt%

ND = Not Detected

Gasoline Range: C4 - C13	Indicators: Isooctane, Olefins
Diesel Range: C7 - C20	Indicators: Pristane, Phytane
Condensate Range: C2 - C25 +	Indicators: No Olefins, Light & Heavies
Heavy Oil: C20 +	

Comments: Analysis performed on hydrocarbon layer.

Joseph Ponminissery, Laboratory Director

Quality Assurance: The above analyses are performed in accordance with ASTM, UOP, GPA guidelines for quality assurance, unless otherwise stated. The test results apply to the sample as received.

Detailed Hydrocarbon Analysis Report

Operator: Hoang Phung

Raw File: H:\COMPASS\PIANO Dragon II Export\CDF Files Dragon II\1030_HGC26\26040367-001A TopLayer HKP.CDF

Sample File: H:\COMPASS\PIANO Dragon II Export\DHA Results Dragon II\26040367-001A TopLayer HKP_04202026_1417.DHA *

Reference File: H:\COMPASS\PIANO Dragon II Export\Reference DHAs Dragon II\19-01-199-01 GC26 NaphCond 03.26.26 HKP_03272026_1324.DHA

Sample Name: 26040367-001A TopLayer HKP Analyzed: 20-Apr-26, 14:17:59 Acquired: 17-Apr-26 01:22:29 Peaks Processed: 265 Total Area: 489 Normalization Factor: 100%

Details Report

Total Hydrocarbons by Group and Carbon Number, Mass %

C #	Paraffins	Iso Paraffins	Olefins	Naphthenes	Aromatics	Unknowns	Oxygenates	Total
C1	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C2	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C3	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C4	0.079	0.006	0.000	0.000	0.000	0.000	0.000	0.085
C5	3.404	1.108	0.000	0.616	0.000	0.000	0.000	5.128
C6	12.679	9.947	0.000	9.969	0.937	0.000	0.000	33.532
C7	9.634	10.439	0.001	15.116	3.136	0.000	0.000	38.326
C8	3.061	8.100	0.210	3.405	1.817	0.000	0.000	16.593
C9	0.618	3.177	0.052	0.945	0.379	0.000	0.000	5.171
C10	0.089	0.672	0.005	0.150	0.104	0.000	0.000	1.020
C11	0.015	0.082	0.000	0.000	0.007	0.000	0.000	0.104
C12	0.003	0.017	0.000	0.002	0.002	0.000	0.000	0.024
C13	0.001	0.001	0.000	0.000	0.001	0.000	0.000	0.003
C14	0.001	0.000	0.000	0.000	0.000	0.000	0.000	0.001
C15	0.000	0.008	0.000	0.000	0.000	0.000	0.000	0.008
C16	0.001	0.000	0.000	0.000	0.000	0.000	0.000	0.001
C17	0.001	0.000	0.000	0.000	0.000	0.000	0.000	0.001
C18	0.001	0.001	0.000	0.000	0.000	0.000	0.000	0.002
C19	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C20	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C21	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C22	0.000	0.001	0.000	0.000	0.000	0.000	0.000	0.001
C23	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C24	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C25	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C26	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C27	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C28	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C29	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C30	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
Total	29.587	33.559	0.268	30.203	6.383	0.000	0.000	100.000

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C1	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C2	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C3	0.001	0.000	0.000	0.000	0.000	0.000	0.000	0.001
C4	0.129	0.009	0.000	0.000	0.000	0.000	0.000	0.138
C5	4.492	1.463	0.000	0.836	0.000	0.000	0.000	6.791
C6	14.007	10.989	0.000	11.278	1.143	0.000	0.000	37.417
C7	9.154	9.920	0.001	14.656	3.240	0.000	0.000	36.971
C8	2.551	6.752	0.177	2.888	1.631	0.000	0.000	13.999
C9	0.459	2.375	0.040	0.712	0.299	0.000	0.000	3.885
C10	0.060	0.469	0.004	0.103	0.073	0.000	0.000	0.709
C11	0.009	0.053	0.000	0.000	0.005	0.000	0.000	0.067
C12	0.002	0.012	0.000	0.001	0.002	0.000	0.000	0.017
C13	0.000	0.001	0.000	0.000	0.000	0.000	0.000	0.001
C14	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C15	0.000	0.004	0.000	0.000	0.000	0.000	0.000	0.004
C16	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C17	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C18	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C19	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C20	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C21	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C22	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C23	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C24	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C25	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C26	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C27	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C28	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C29	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
C30	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
Total	30.864	32.047	0.222	30.474	6.393	0.000	0.000	100.000

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Individual Component Results

Component	Group	Retention Time	Retention Index	Area	Mass %	Vol %	Mole %	BP (F)	BP (C)	CAS #
propane	P3	8.599	300.00	0.0	0.000	0.000	0.001	(43.7)	(42.0)	74-98-6
i-butane	I4	9.070	355.11	0.0	0.006	0.007	0.009	10.9	(11.7)	75-28-5
n-butane	P4	9.473	400.00	0.4	0.079	0.098	0.129	31.1	(0.5)	106-97-8
2,2-dimethylpropane	I5	9.676	409.87	0.0	0.000	0.001	0.001	49.1	9.5	463-82-1
i-pentane	I5	10.951	467.37	5.0	1.108	1.274	1.462	82.1	27.8	78-78-4
n-pentane	P5	11.748	500.00	15.4	3.404	3.875	4.492	96.9	36.1	109-66-0
2,2-dimethylbutane	I6	13.274	530.11	0.4	0.082	0.091	0.090	121.5	49.7	75-83-2
cyclopentane	N5	14.863	557.96	2.9	0.616	0.589	0.836	120.7	49.3	287-92-3
2,3-dimethylbutane	I6	14.961	559.58	2.9	0.622	0.670	0.687	136.4	58.0	79-29-8
2-methylpentane	I6	15.215	563.73	26.6	5.726	6.231	6.326	140.5	60.3	107-83-5
3-methylpentane	I6	16.233	579.68	16.4	3.517	3.774	3.886	145.9	63.3	96-14-0
n-hexane	P6	17.628	600.00	60.4	12.679	13.671	14.007	155.7	68.7	110-54-3
C7-Olefin	O7	19.684	621.59	0.0	0.001	0.001	0.001	198.9	92.7	
2,2-dimethylpentane	I7	19.887	623.60	1.2	0.245	0.251	0.233	174.5	79.2	590-35-2
methylcyclopentane	N6	20.108	625.77	25.5	5.175	4.928	5.854	161.2	71.8	96-37-7
2,4-dimethylpentane	I7	20.504	629.58	3.7	0.770	0.794	0.732	176.9	80.5	108-08-7
2,2,3-trimethylbutane	I7	21.098	635.17	0.2	0.035	0.036	0.033	177.6	80.9	464-06-2
benzene	A6	22.622	648.81	5.4	0.937	0.763	1.143	176.2	80.1	71-43-2
3,3-dimethylpentane	I7	23.322	654.77	0.6	0.120	0.123	0.114	186.9	86.1	562-49-2
cyclohexane	N6	23.725	658.13	23.6	4.794	4.390	5.424	177.3	80.7	110-82-7
2-methylhexane	I7	24.873	667.37	18.6	3.846	3.956	3.655	194.1	90.1	591-76-4
2,3-dimethylpentane	I7	25.074	668.95	6.0	1.238	1.277	1.176	193.6	89.8	565-59-3
1,1-dimethylcyclopentane	N7	25.410	671.55	5.0	1.015	0.937	0.984	189.5	87.5	1638-26-2
3-methylhexane	I7	25.987	675.95	19.0	3.922	4.035	3.727	197.3	91.9	589-34-4
1c,3-dimethylcyclopentane	N7	26.796	681.95	7.0	1.413	1.322	1.370	195.4	90.8	2532-58-3
1t,3-dimethylcyclopentane	N7	27.169	684.65	6.4	1.291	1.208	1.252	197.1	91.7	1759-58-6
3-ethylpentane	I7	27.353	685.98	1.3	0.263	0.270	0.250	200.2	93.5	617-78-7
1t,2-dimethylcyclopentane	N7	27.529	687.23	9.7	1.958	1.831	1.898	197.4	91.9	822-50-4
2,2,4-trimethylpentane	I8	27.786	689.04	0.1	0.019	0.020	0.016	210.6	99.2	540-84-1
n-heptane	P7	29.386	700.00	47.0	9.634	10.107	9.154	209.2	98.4	142-82-5
1c,2-dimethylcyclopentane	N7	32.103	719.79	1.1	0.217	0.203	0.210	211.2	99.5	1192-18-3
methylcyclohexane	N7	32.217	720.58	44.4	8.918	8.262	8.647	213.7	100.9	108-87-2
2,2-dimethylhexane	I8	32.678	723.76	3.9	0.792	0.795	0.661	224.3	106.8	590-73-8
ethylcyclopentane	N7	33.973	732.45	1.5	0.304	0.279	0.295	218.2	103.5	1640-89-7
2,5-dimethylhexane	I8	34.198	733.93	1.5	0.308	0.310	0.257	228.4	109.1	592-13-2
2,2,3-trimethylpentane	I8	34.361	735.00	0.2	0.032	0.032	0.026	229.7	109.9	564-02-3
2,4-dimethylhexane	I8	34.488	735.83	2.4	0.493	0.497	0.411	229.0	109.4	589-43-5
1c,2t,4-trimethylcyclopentane	N8	35.338	741.27	3.1	0.618	0.585	0.525	242.1	116.7	4850-28-6
3,3-dimethylhexane	I8	35.581	742.81	0.5	0.094	0.094	0.078	233.5	112.0	563-16-6
C8-Olefin	O8	36.268	747.08	0.0	0.003	0.003	0.002	248.0	120.0	
1t,2c,3-trimethylcyclopentane	N8	36.487	748.43	2.4	0.473	0.447	0.401	230.7	110.4	15890-40-1

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Individual Component Results

Component	Group	Retention Time	Retention Index	Area	Mass %	Vol %	Mole %	BP (F)	BP (C)	CAS #
2,3,4-trimethylpentane	I8	36.964	751.34	0.3	0.071	0.072	0.059	236.2	113.5	565-75-3
toluene	A7	37.503	754.58	17.4	3.136	2.592	3.240	231.1	110.6	108-88-3
2,3-dimethylhexane	I8	38.758	761.94	2.2	0.455	0.459	0.379	240.1	115.6	584-94-1
2-methyl-3-ethylpentane	I8	38.944	763.02	0.3	0.065	0.066	0.055	240.1	115.6	609-26-7
C8-Olefin	O8	39.348	765.33	0.0	0.003	0.003	0.002	248.0	120.0	
2-methylheptane	I8	39.753	767.61	10.3	2.111	2.122	1.759	243.8	117.7	592-27-8
4-methylheptane	I8	39.993	768.96	2.8	0.577	0.580	0.481	243.9	117.7	589-53-7
3-methyl-3-ethylpentane	I8	40.173	769.96	0.3	0.054	0.054	0.045	240.1	115.6	1067-08-9
3,4-dimethylhexane	I8	40.263	770.47	0.4	0.081	0.082	0.068	243.9	117.7	583-48-2
1c,2t,3-trimethylcyclopentane	N8	40.643	772.57	0.3	0.066	0.062	0.056	243.5	117.5	19374-46-0
3-methylheptane	I8	41.003	774.55	5.7	1.172	1.178	0.977	246.1	118.9	589-81-1
3-ethylhexane	I8	41.203	775.63	8.7	1.776	1.786	1.480	245.4	118.5	619-99-8
C8-Olefin	O8	41.330	776.32	0.4	0.078	0.073	0.066	248.0	120.0	
1t,4-dimethylcyclohexane	N8	41.534	777.42	3.5	0.700	0.654	0.594	246.8	119.4	2207-04-7
1,1-dimethylcyclohexane	N8	42.495	782.54	1.3	0.254	0.235	0.215	247.2	119.6	590-66-9
2,2,5-trimethylhexane	I9	42.933	784.83	0.1	0.013	0.013	0.010	255.4	124.1	3522-94-9
3c-ethylmethylcyclopentane	N8	43.160	786.02	0.4	0.072	0.066	0.061	250.0	121.1	2613-65-2
3t-ethylmethylcyclopentane	N8	43.553	788.04	0.3	0.064	0.059	0.054	250.0	121.1	2613-66-3
octene-1	O8	43.782	789.22	0.6	0.125	0.124	0.106	256.1	124.5	111-66-0
1,1-methylethylcyclopentane	N8	44.178	791.23	0.3	0.051	0.047	0.043	250.8	121.5	16747-50-5
1t,2-dimethylcyclohexane	N8	44.678	793.75	3.3	0.660	0.614	0.560	254.2	123.4	6876-23-9
n-octane	P8	45.943	800.00	14.9	3.061	3.123	2.551	258.2	125.7	111-65-9
1c,4-dimethylcyclohexane	N8	46.090	800.75	1.6	0.312	0.291	0.265	255.8	124.3	624-29-3
C9-IsoParaffin	I9	47.544	808.02	0.1	0.013	0.012	0.010	297.6	147.6	
C9-IsoParaffin	I9	47.618	808.38	0.2	0.036	0.034	0.028	297.6	147.6	
c-octene-2	O8	48.119	810.84	0.0	0.001	0.001	0.001	32.0	0.0	7642-04-8
C9-IsoParaffin	I9	48.376	812.09	0.0	0.003	0.003	0.002	297.6	147.6	
2,3,5-trimethylhexane	I9	48.848	814.36	0.1	0.026	0.026	0.019	268.4	131.4	1069-53-0
N2	N8	49.343	816.72	0.1	0.023	0.021	0.019	267.2	130.7	
C9-IsoParaffin	I9	49.716	818.49	0.1	0.012	0.011	0.009	297.6	147.6	
1c,2-dimethylcyclohexane	N8	50.143	820.49	0.3	0.065	0.060	0.055	265.5	129.7	2207-01-4
1,1,4-trimethylcyclohexane	N9	51.051	824.69	1.3	0.257	0.239	0.193	275.0	135.0	7094-27-1
C9-IsoParaffin	I9	51.258	825.64	0.1	0.029	0.027	0.022	297.6	147.6	
C9-IsoParaffin	I9	51.481	826.66	0.0	0.008	0.007	0.006	297.6	147.6	
C9-IsoParaffin	I9	51.723	827.75	0.0	0.007	0.006	0.005	297.6	147.6	
2,2,3-trimethylhexane	I9	52.239	830.08	4.4	0.891	0.880	0.661	271.2	132.9	16747-25-4
C9-IsoParaffin	I9	52.617	831.77	1.3	0.254	0.234	0.193	297.6	147.6	
C9-IsoParaffin	I9	53.066	833.76	0.1	0.014	0.012	0.010	297.6	147.6	
4,4-dimethylheptane	I9	53.553	835.90	1.5	0.307	0.306	0.228	271.2	132.9	1068-19-5
3,5-dimethylheptane	I9	54.211	838.76	1.2	0.232	0.242	0.175	278.6	137.0	4032-86-4
n-propylcyclopentane	N8	54.611	840.48	0.2	0.047	0.043	0.040	267.7	131.0	2040-96-2

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Sample File: H:\COMPASS\PIANO Dragon II Export\DHA Results Dragon II\26040367-001A TopLayer HKP_04202026_1417.DHA *

Reference File: H:\COMPASS\PIANO Dragon II Export\Reference DHAs Dragon II\19-01-199-01 GC26 NaphCond 03.26.26 HKP_03272026_1324.DHA

Sample Name: 26040367-001A TopLayer HKP Analyzed: 20-Apr-26, 14:17:59 Acquired: 17-Apr-26 01:22:29 Peaks Processed: 265 Total Area: 489 Normalization Factor: 100%

Individual Component Results

Component	Group	Retention Time	Retention Index	Area	Mass %	Vol %	Mole %	BP (F)	BP (C)	CAS #
C9-IsoParaffin	I9	54.850	841.51	0.1	0.016	0.015	0.012	297.6	147.6	
C9-IsoParaffin	I9	55.173	842.88	0.2	0.030	0.028	0.023	297.6	147.6	
C9-IsoParaffin	I9	55.279	843.33	0.1	0.011	0.010	0.009	297.6	147.6	
C9-IsoParaffin	I9	55.576	844.59	0.1	0.016	0.015	0.012	297.6	147.6	
ethylbenzene	A8	56.700	849.28	1.2	0.212	0.175	0.190	277.2	136.2	100-41-4
C9-IsoParaffin	I9	57.011	850.56	0.2	0.037	0.034	0.028	297.6	147.6	
C9-IsoParaffin	I9	57.140	851.09	0.1	0.023	0.022	0.018	297.6	147.6	
1c,2t,4t-trimethylcyclohexane	N9	57.463	852.41	0.9	0.175	0.165	0.132	32.0	0.0	7667-60-9
C9-IsoParaffin	I9	58.008	854.62	0.0	0.003	0.003	0.003	297.6	147.6	
C9-IsoParaffin	I9	58.293	855.77	0.0	0.002	0.002	0.001	297.6	147.6	
C9-IsoParaffin	I9	58.574	856.89	0.1	0.011	0.010	0.009	297.6	147.6	
m-xylene	A8	59.081	858.91	5.1	0.935	0.766	0.839	282.4	139.1	108-38-3
p-xylene	A8	59.393	860.14	1.9	0.334	0.277	0.300	281.0	138.4	106-42-3
2,3-dimethylheptane	I9	59.753	861.56	1.5	0.307	0.303	0.228	284.9	140.5	3074-71-3
3,4-dimethylheptane	I9	60.313	863.75	0.1	0.019	0.019	0.014	285.1	140.6	922-28-1
C9-IsoParaffin	I9	60.608	864.89	0.4	0.082	0.076	0.062	297.6	147.6	
C9-IsoParaffin	I9	61.166	867.04	0.1	0.027	0.025	0.021	297.6	147.6	
C9-IsoParaffin	I9	61.414	867.99	0.0	0.002	0.002	0.001	297.6	147.6	
4-methyloctane	I9	62.020	870.28	0.9	0.181	0.179	0.134	288.4	142.4	2216-34-4
2-methyloctane	I9	62.307	871.37	1.0	0.206	0.203	0.153	289.9	143.3	3221-61-2
C9-IsoParaffin	I9	62.873	873.48	0.1	0.022	0.021	0.017	297.6	147.6	
3-ethylheptane	I9	63.628	876.28	0.3	0.067	0.067	0.050	289.4	143.0	15869-80-4
3-methyloctane	I9	64.003	877.66	1.1	0.228	0.226	0.170	291.6	144.2	2216-33-3
C9-IsoParaffin	I9	64.511	879.51	0.0	0.005	0.005	0.004	297.6	147.6	
C9-IsoParaffin	I9	64.712	880.24	0.1	0.012	0.011	0.009	297.6	147.6	
o-xylene	A8	64.961	881.14	1.9	0.336	0.273	0.302	292.0	144.4	95-47-6
C9-IsoParaffin	I9	65.392	882.69	0.1	0.014	0.013	0.011	297.6	147.6	
C9-IsoParaffin	I9	66.005	884.87	0.0	0.008	0.008	0.006	297.6	147.6	
N18	N9	66.541	886.77	0.2	0.046	0.042	0.035	297.9	147.7	
N19	N9	66.756	887.52	0.9	0.182	0.166	0.137	298.1	147.8	
N20	N9	67.254	889.26	0.6	0.126	0.115	0.095	299.1	148.4	
1-Nonene	O9	67.913	891.55	0.0	0.006	0.006	0.005	274.1	134.5	124-11-8
i-butylcyclopentane	N9	68.247	892.70	0.1	0.012	0.011	0.009	298.3	148.0	3788-32-7
N22	N9	68.550	893.73	0.0	0.003	0.003	0.002	301.6	149.8	
t-nonene-3	O9	69.376	896.54	0.1	0.018	0.018	0.014	32.0	0.0	20063-92-7
C9-IsoParaffin	I9	69.791	897.94	0.0	0.003	0.003	0.002	297.6	147.6	
cis-3-Nonene	O9	69.948	898.46	0.0	0.009	0.009	0.007	297.9	147.7	20237-46-1
n-nonane	P9	70.408	900.00	3.1	0.618	0.614	0.459	303.5	150.8	111-84-2
1,1-methylethylcyclohexane	N9	70.816	902.72	0.5	0.095	0.087	0.072	305.9	152.2	4926-90-3
t-nonene-2	O9	71.306	905.98	0.1	0.019	0.018	0.014	32.0	0.0	6434-78-2
N25	N9	71.428	906.79	0.1	0.013	0.012	0.010	305.7	152.1	

Detailed Hydrocarbon Analysis Report

Operator: Hoang Phung

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Reference File: H:\COMPASS\PIANO Dragon II Export\Reference DHAs Dragon II\19-01-199-01 GC26 NaphCond 03.26.26 HKP_03272026_1324.DHA

Sample Name: 26040367-001A TopLayer HKP Analyzed: 20-Apr-26, 14:17:59 Acquired: 17-Apr-26 01:22:29 Peaks Processed: 265 Total Area: 489 Normalization Factor: 100%

Individual Component Results

Component	Group	Retention Time	Retention Index	Area	Mass %	Vol %	Mole %	BP (F)	BP (C)	CAS #
C10-IsoParaffin	I10	71.718	908.69	0.0	0.002	0.002	0.001	340.2	171.2	
i-propylbenzene	A9	72.294	912.47	0.1	0.024	0.020	0.019	306.3	152.4	98-82-8
C10-IsoParaffin	I10	72.861	916.15	0.3	0.064	0.055	0.046	340.2	171.2	
i-propylcyclohexane	N9	73.039	917.30	0.1	0.010	0.009	0.007	310.6	154.8	696-29-7
C10-IsoParaffin	I10	73.316	919.09	0.2	0.031	0.027	0.022	340.2	171.2	
C10-IsoParaffin	I10	73.768	921.99	0.1	0.014	0.012	0.010	340.2	171.2	
C10-IsoParaffin	I10	74.120	924.24	0.1	0.021	0.018	0.015	340.2	171.2	
2,2-dimethyloctane	I10	74.184	924.64	0.2	0.034	0.033	0.023	314.4	156.9	15869-87-1
N28	N9	74.528	926.82	0.1	0.012	0.011	0.009	314.4	156.9	
N29	N9	74.948	929.47	0.1	0.014	0.013	0.011	315.1	157.3	
C10-IsoParaffin	I10	75.343	931.95	0.6	0.119	0.101	0.085	340.2	171.2	
I12	I10	75.511	933.01	0.2	0.036	0.035	0.024	311.4	155.2	
C10-IsoParaffin	I10	75.820	934.93	0.0	0.002	0.002	0.001	340.2	171.2	
C10-IsoParaffin	I10	76.133	936.87	0.1	0.014	0.012	0.010	340.2	171.2	
2,6-dimethyloctane	I10	76.257	937.64	0.0	0.008	0.008	0.006	320.7	160.4	2051-30-1
C10-IsoParaffin	I10	76.429	938.71	0.1	0.015	0.013	0.011	340.2	171.2	
C10-IsoParaffin	I10	76.780	940.87	0.1	0.011	0.009	0.008	340.2	171.2	
3,3-dimethyloctane	I10	77.063	942.61	0.4	0.072	0.070	0.048	322.2	161.2	4110-44-5
C10-IsoParaffin	I10	77.374	944.51	0.1	0.017	0.014	0.012	340.2	171.2	
n-propylbenzene	A9	77.692	946.44	0.3	0.049	0.040	0.038	318.6	159.2	103-65-1
3-methyl-5-ethylheptane	I10	78.024	948.45	0.1	0.015	0.014	0.010	316.8	158.2	52896-90-9
N32	N10	78.259	949.87	0.2	0.042	0.038	0.029	324.9	162.7	
C10-IsoParaffin	I10	78.661	952.29	0.1	0.011	0.009	0.008	340.2	171.2	
1,3-methylethylbenzene	A9	78.994	954.28	0.4	0.077	0.063	0.061	322.4	161.3	620-14-4
1,4-methylethylbenzene	A9	79.316	956.20	0.2	0.034	0.028	0.027	323.6	162.0	622-96-8
C10-IsoParaffin	I10	79.551	957.60	0.0	0.001	0.001	0.001	340.2	171.2	
N33	N10	79.921	959.78	0.1	0.016	0.014	0.011	327.8	164.3	
1,3,5-trimethylbenzene	A9	80.268	961.83	0.3	0.058	0.048	0.046	328.5	164.7	108-67-8
2,3-dimethyloctane	I10	80.409	962.66	0.1	0.022	0.022	0.015	327.8	164.3	7146-60-3
N34	N10	80.586	963.69	0.1	0.011	0.010	0.007	330.6	165.9	
C10-IsoParaffin	I10	80.632	963.96	0.0	0.008	0.007	0.006	340.2	171.2	
C10-IsoParaffin	I10	80.963	965.90	0.0	0.009	0.008	0.007	340.2	171.2	
C10-IsoParaffin	I10	81.084	966.60	0.0	0.003	0.003	0.002	340.2	171.2	
5-methylnonane	I10	81.323	967.99	0.1	0.011	0.011	0.008	329.2	165.1	15869-85-9
C10-IsoParaffin	I10	81.583	969.50	0.2	0.035	0.030	0.025	340.2	171.2	
1,2-methylethylbenzene	A9	81.729	970.34	0.1	0.017	0.014	0.013	329.3	165.2	611-14-3
2-methylnonane	I10	81.998	971.89	0.2	0.032	0.031	0.021	332.7	167.0	871-83-0
C10-IsoParaffin	I10	82.274	973.48	0.0	0.004	0.004	0.003	340.2	171.2	
3-ethyloctane	I10	82.471	974.60	0.0	0.009	0.009	0.006	331.7	166.5	5881-17-4
N35	N10	82.699	975.91	0.1	0.012	0.011	0.008	335.3	168.5	
3-methylnonane	I10	82.966	977.42	0.1	0.026	0.025	0.017	334.0	167.8	5911-04-6

Detailed Hydrocarbon Analysis Report

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Sample Name: 26040367-001A TopLayer HKP Analyzed: 20-Apr-26, 14:17:59 Acquired: 17-Apr-26 01:22:29 Peaks Processed: 265 Total Area: 489 Normalization Factor: 100%

Individual Component Results

Component	Group	Retention Time	Retention Index	Area	Mass %	Vol %	Mole %	BP (F)	BP (C)	CAS #
C10-IsoParaffin	I10	83.171	978.59	0.0	0.009	0.007	0.006	340.2	171.2	
I18	I10	83.590	980.96	0.0	0.003	0.003	0.002	337.0	169.4	
1,2,4-Trimethylbenzene	A9	84.018	983.37	0.6	0.101	0.083	0.080	336.9	169.4	95-63-6
tert-Butylcyclohexane	N10	84.301	984.96	0.1	0.023	0.020	0.016	332.6	167.0	3178-22-1
i-butylcyclohexane	N10	84.694	987.15	0.1	0.023	0.021	0.016	340.3	171.3	1678-98-4
2-Methyl-1-nonene	O10	84.861	988.08	0.0	0.002	0.002	0.002	212.0	100.0	2980-71-4
I23	I10	85.041	989.08	0.0	0.004	0.004	0.003	341.0	171.7	
N37	N10	85.234	990.15	0.0	0.004	0.004	0.003	341.4	171.9	
decene-1	O10	85.411	991.13	0.0	0.003	0.003	0.002	339.1	170.6	872-05-9
C10-IsoParaffin	I10	85.978	994.25	0.0	0.006	0.005	0.004	340.2	171.2	
i-butylbenzene	A10	86.198	995.45	0.1	0.011	0.009	0.008	343.0	172.8	538-93-2
sec-butylbenzene	A10	86.673	998.04	0.1	0.009	0.008	0.007	344.0	173.3	135-98-8
C10-IsoParaffin	I10	86.786	998.66	0.0	0.004	0.003	0.003	340.2	171.2	
n-decane	P10	87.033	1000.00	0.5	0.089	0.087	0.060	345.5	174.2	124-18-5
C11-IsoParaffin	I11	87.305	1002.44	0.0	0.004	0.003	0.003	378.5	192.5	
C11-IsoParaffin	I11	87.625	1005.29	0.0	0.003	0.002	0.002	378.5	192.5	
1,2,3-trimethylbenzene	A9	87.829	1007.11	0.1	0.015	0.012	0.012	349.0	176.1	526-73-8
1,3-methyl-i-propylbenzene	A10	88.118	1009.67	0.0	0.009	0.007	0.006	347.1	175.1	535-77-3
1,4-methyl-i-propylbenzene	A10	88.532	1013.33	0.0	0.004	0.004	0.003	350.8	177.1	99-87-6
C11-IsoParaffin	I11	88.758	1015.32	0.0	0.002	0.002	0.001	378.5	192.5	
C11-IsoParaffin	I11	89.048	1017.86	0.0	0.002	0.002	0.002	378.5	192.5	
2-3-dihydroindene	A9	89.168	1018.91	0.0	0.004	0.003	0.003	352.1	177.9	496-11-7
C11-IsoParaffin	I11	89.382	1020.79	0.0	0.004	0.003	0.003	378.5	192.5	
C11-IsoParaffin	I11	89.643	1023.06	0.0	0.003	0.002	0.002	378.5	192.5	
1,2-methyl-i-propylbenzene	A10	89.864	1024.99	0.0	0.009	0.007	0.006	352.7	178.2	1074-17-5
C11-IsoParaffin	I11	90.221	1028.08	0.1	0.016	0.013	0.011	378.5	192.5	
C11-IsoParaffin	I11	90.618	1031.50	0.0	0.002	0.001	0.001	378.5	192.5	
n-Butylcyclohexane	N10	90.833	1033.36	0.1	0.019	0.017	0.013	358.1	181.1	1678-93-9
C11-IsoParaffin	I11	91.344	1037.73	0.0	0.005	0.004	0.003	378.5	192.5	
1,3-diethylbenzene	A10	91.577	1039.72	0.0	0.004	0.004	0.003	358.1	181.1	141-93-5
1,3-methyl-n-propylbenzene	A10	91.909	1042.55	0.1	0.010	0.008	0.007	359.6	182.0	1074-43-7
C11-IsoParaffin	I11	92.084	1044.03	0.0	0.002	0.001	0.001	378.5	192.5	
C11-IsoParaffin	I11	92.189	1044.92	0.0	0.004	0.003	0.002	378.5	192.5	
1,4-methyl-n-propylbenzene	A10	92.376	1046.50	0.0	0.004	0.003	0.003	362.2	183.4	1074-55-1
n-butylbenzene	A10	92.508	1047.62	0.0	0.003	0.003	0.002	361.9	183.3	104-51-8
1,3-dimethyl-5-ethylbenzene	A10	92.743	1049.59	0.0	0.006	0.005	0.004	362.5	183.6	934-74-7
C11-IsoParaffin	I11	92.993	1051.70	0.0	0.002	0.001	0.001	378.5	192.5	
C11-IsoParaffin	I11	93.245	1053.81	0.0	0.002	0.002	0.001	378.5	192.5	
C11-IsoParaffin	I11	93.412	1055.20	0.0	0.006	0.005	0.004	378.5	192.5	
1,2-methyl-n-propylbenzene	A10	93.740	1057.94	0.0	0.006	0.005	0.004	364.9	185.0	1074-17-5
C11-IsoParaffin	I11	93.927	1059.49	0.0	0.002	0.001	0.001	378.5	192.5	

Detailed Hydrocarbon Analysis Report

Operator: Hoang Phung

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Sample Name: 26040367-001A TopLayer HKP Analyzed: 20-Apr-26, 14:17:59 Acquired: 17-Apr-26 01:22:29 Peaks Processed: 265 Total Area: 489 Normalization Factor: 100%

Individual Component Results

Component	Group	Retention Time	Retention Index	Area	Mass %	Vol %	Mole %	BP (F)	BP (C)	CAS #
C11-IsoParaffin	I11	94.169	1061.50	0.0	0.001	0.001	0.001	378.5	192.5	
5-methyldecane	I11	94.328	1062.82	0.0	0.004	0.004	0.002	367.0	186.1	13151-35-4
C11-IsoParaffin	I11	94.687	1065.78	0.0	0.004	0.003	0.003	378.5	192.5	
1,4-dimethyl-2-ethylbenzene	A10	94.941	1067.87	0.0	0.003	0.003	0.002	368.4	186.9	1758-88-9
1,3-dimethyl-4-ethylbenzene	A10	95.108	1069.25	0.0	0.006	0.005	0.004	370.8	188.2	874-41-9
C11-IsoParaffin	I11	95.358	1071.29	0.0	0.001	0.001	0.001	378.5	192.5	
1,2-dimethyl-4-ethylbenzene	A10	95.829	1075.14	0.0	0.007	0.006	0.005	373.1	189.5	934-80-5
C11-IsoParaffin	I11	95.998	1076.51	0.0	0.001	0.001	0.001	378.5	192.5	
C11-IsoParaffin	I11	96.350	1079.37	0.0	0.002	0.002	0.001	378.5	192.5	
C11-IsoParaffin	I11	96.457	1080.23	0.0	0.002	0.002	0.001	378.5	192.5	
C11-IsoParaffin	I11	96.854	1083.44	0.0	0.001	0.001	0.001	378.5	192.5	
C11-IsoParaffin	I11	97.178	1086.05	0.0	0.002	0.002	0.001	378.5	192.5	
C11-IsoParaffin	I11	97.322	1087.20	0.0	0.002	0.002	0.001	378.5	192.5	
C11-IsoParaffin	I11	97.639	1089.74	0.0	0.002	0.002	0.001	378.5	192.5	
1,4-methyl-t-butylbenzene	A11	97.854	1091.46	0.0	0.002	0.002	0.002	32.0	0.0	98-51-1
1,2-Dimethyl-3-ethylbenzene	A10	98.097	1093.39	0.0	0.003	0.002	0.002	381.1	194.0	933-98-2
C11-IsoParaffin	I11	98.278	1094.83	0.0	0.001	0.001	0.001	378.5	192.5	
1,2-ethyl-i-propylbenzene	A11	98.533	1096.85	0.0	0.003	0.002	0.002	32.0	0.0	16021-20-8
n-undecane	P11	98.931	1100.00	0.1	0.015	0.015	0.009	384.6	195.9	1120-21-4
C12-IsoParaffin	I12	99.264	1103.54	0.0	0.001	0.001	0.001	457.3	236.3	
1,2,4,5-tetramethylbenzene	A10	99.403	1105.01	0.0	0.002	0.002	0.002	386.3	196.8	95-93-2
1,2,3,5-tetramethylbenzene	A10	99.783	1109.02	0.0	0.002	0.002	0.001	388.5	198.0	527-53-7
C12-IsoParaffin	I12	99.990	1111.21	0.0	0.003	0.002	0.002	457.3	236.3	
C12-IsoParaffin	I12	100.639	1118.02	0.0	0.001	0.001	0.001	457.3	236.3	
1,2-methyl-t-butylbenzene	A11	100.930	1121.06	0.0	0.002	0.001	0.001	32.0	0.0	1074-92-6
5-methylindan	A10	101.517	1127.16	0.0	0.001	0.001	0.001	32.0	0.0	874-35-1
C12-IsoParaffin	I12	101.713	1129.20	0.0	0.002	0.001	0.001	457.3	236.3	
C12-IsoParaffin	I12	101.908	1131.21	0.0	0.001	0.001	0.001	457.3	236.3	
2-methylindan	A10	102.443	1136.72	0.0	0.003	0.003	0.002	368.6	187.0	824-63-5
C12-IsoParaffin	I12	102.791	1140.29	0.0	0.001	0.001	0.000	457.3	236.3	
1,3-di-i-propylbenzene	A12	103.073	1143.18	0.0	0.001	0.001	0.001	397.8	203.2	3748-13-8
C12-IsoParaffin	I12	103.581	1148.35	0.0	0.002	0.001	0.001	457.3	236.3	
1t-M-2-(4-MP)cyclopentane	N12	103.990	1152.50	0.0	0.002	0.002	0.001	32.0	0.0	66553-50-2
C12-IsoParaffin	I12	104.453	1157.17	0.0	0.001	0.001	0.001	457.3	236.3	
C12-IsoParaffin	I12	104.950	1162.17	0.0	0.001	0.001	0.001	457.3	236.3	
naphthalene	A10	105.354	1166.22	0.0	0.002	0.001	0.001	424.4	218.0	91-20-3
C12-IsoParaffin	I12	105.628	1168.95	0.0	0.001	0.001	0.001	457.3	236.3	
1-t-butyl-3,5-dimethylbenzene	A12	105.988	1172.53	0.0	0.001	0.001	0.001	32.0	0.0	98-19-1
I45	I12	106.630	1178.89	0.0	0.002	0.001	0.001	413.4	211.9	
I47	I12	107.054	1183.07	0.0	0.001	0.001	0.001	415.2	212.9	
n-dodecane	P12	108.790	1200.00	0.0	0.003	0.003	0.002	421.3	216.3	112-40-3

Detailed Hydrocarbon Analysis Report

Operator: Hoang Phung

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Reference File: H:\COMPASS\PIANO Dragon II Export\Reference DHAs Dragon II\19-01-199-01 GC26 NaphCond 03.26.26 HKP_03272026_1324.DHA

Sample Name: 26040367-001A TopLayer HKP Analyzed: 20-Apr-26, 14:17:59 Acquired: 17-Apr-26 01:22:29 Peaks Processed: 265 Total Area: 489 Normalization Factor: 100%

Individual Component Results

Component	Group	Retention Time	Retention Index	Area	Mass %	Vol %	Mole %	BP (F)	BP (C)	CAS #
C13-IsoParaffin	I13	109.181	1203.69	0.0	0.001	0.001	0.001	478.6	248.1	
1,3,5-triethylbenzene	A12	110.400	1215.12	0.0	0.000	0.000	0.000	420.8	216.0	102-25-0
1,4-methyl-n-pentylbenzene	A12	112.978	1238.86	0.0	0.000	0.000	0.000	32.0	0.0	1595-09-1
1,3-dimethylindane	A11	114.511	1252.73	0.0	0.000	0.000	0.000	407.7	208.7	4175-53-5
1,2,3,4,5-pentamethylbenzene	A11	116.945	1274.37	0.0	0.000	0.000	0.000	449.6	232.0	700-12-9
2-methylnaphthalene	A11	117.572	1279.87	0.0	0.000	0.000	0.000	465.9	241.1	91-57-6
n-tridecane	P13	119.894	1300.00	0.0	0.001	0.001	0.000	455.7	235.4	629-50-5
n-tetradecane	P14	130.188	1400.00	0.0	0.001	0.001	0.000	485.6	252.0	629-59-4
Naphthalene, 1-propyl-	A13	136.950	1490.75	0.0	0.001	0.000	0.000	527.5	275.3	2765-18-6
C15-IsoParaffin	I15	137.611	1499.38	0.0	0.008	0.006	0.004	551.0	288.0	
n-hexadecane	P16	143.499	1600.00	0.0	0.001	0.001	0.000	547.9	286.6	544-76-3
n-heptadecane	P17	148.399	1700.00	0.0	0.001	0.001	0.000	575.6	302.0	629-78-7
C18-IsoParaffin	I18	150.120	1740.83	0.0	0.001	0.001	0.000	602.6	317.0	
n-octadecane	P18	152.649	1800.00	0.0	0.001	0.001	0.000	602.6	317.0	593-45-3
n-nonadecane	P19	156.459	1900.00	0.0	0.000	0.000	0.000	626.0	330.0	629-92-5
n-icosane	P20	159.934	2000.00	0.0	0.000	0.000	0.000	648.9	342.7	112-95-8
n-heneicosane	P21	163.183	2100.00	0.0	0.000	0.000	0.000	673.7	356.5	629-94-7
C22-IsoParaffin	I22	163.941	2123.80	0.0	0.001	0.001	0.000	695.5	368.0	
n-docosane	P22	166.390	2200.00	0.0	0.000	0.000	0.000	694.4	368.0	629-97-0
				489	100.0	100.0	100.0			

Detailed Hydrocarbon Analysis Report

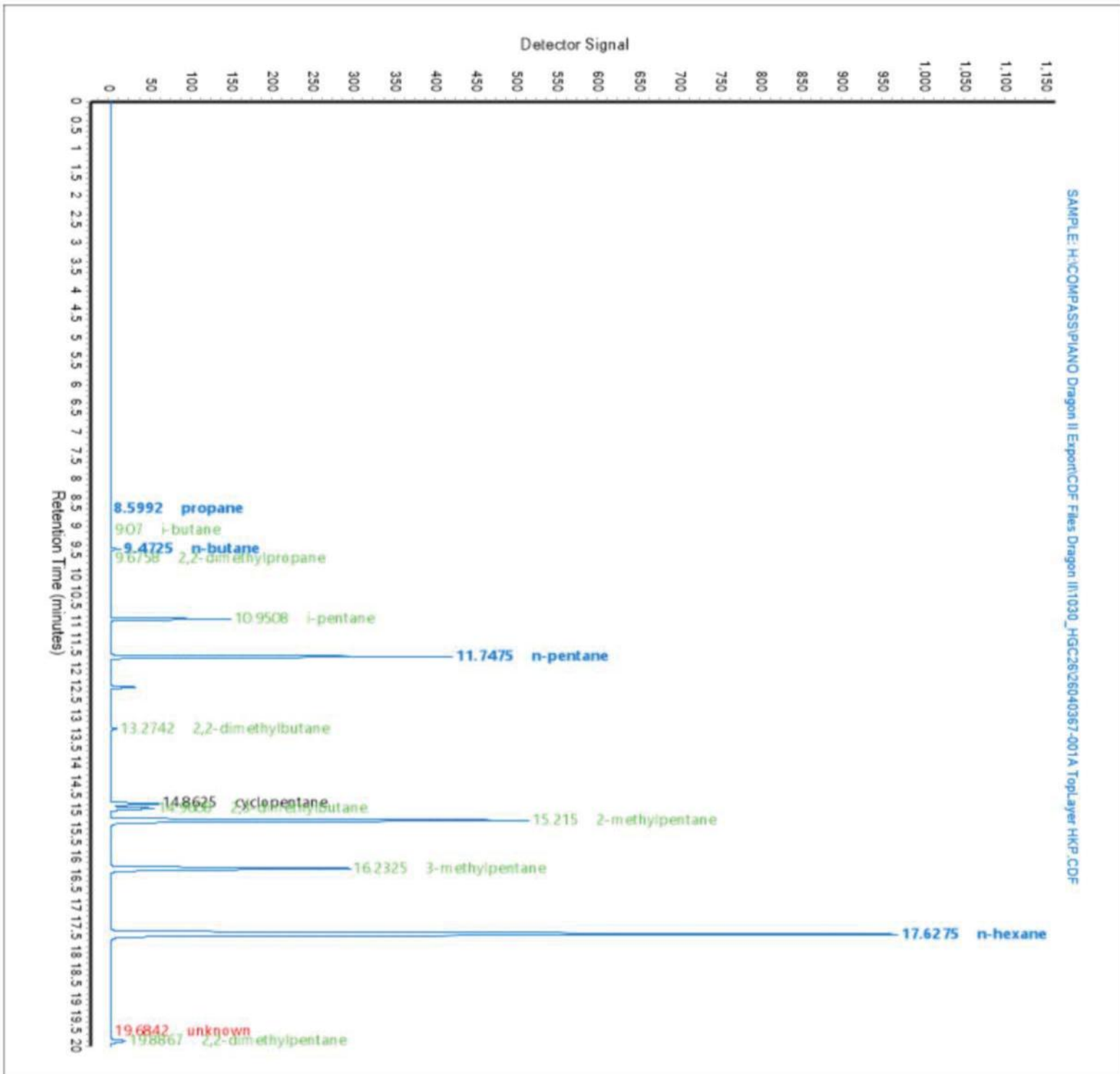
Operator: Hoang Phung

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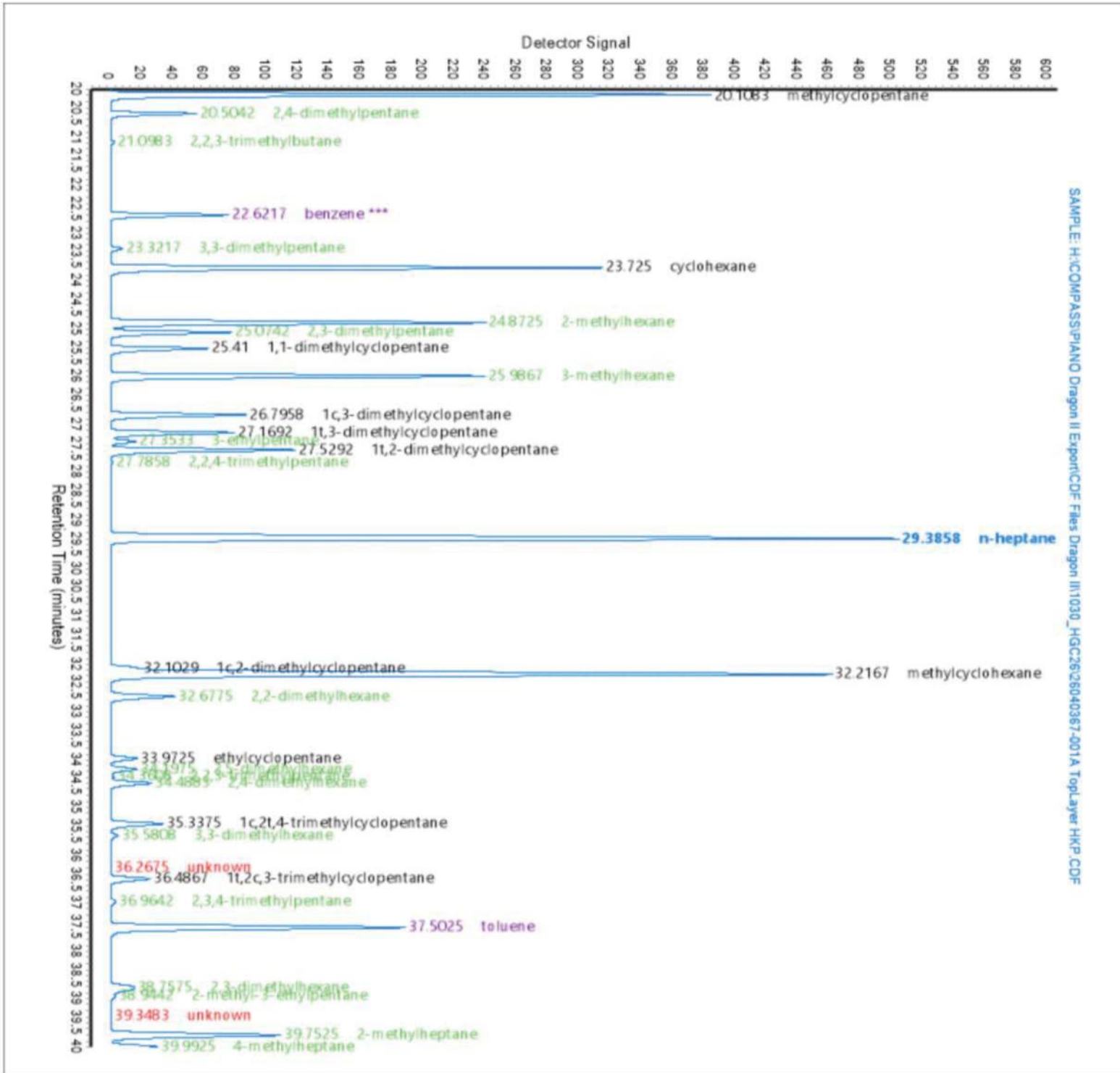
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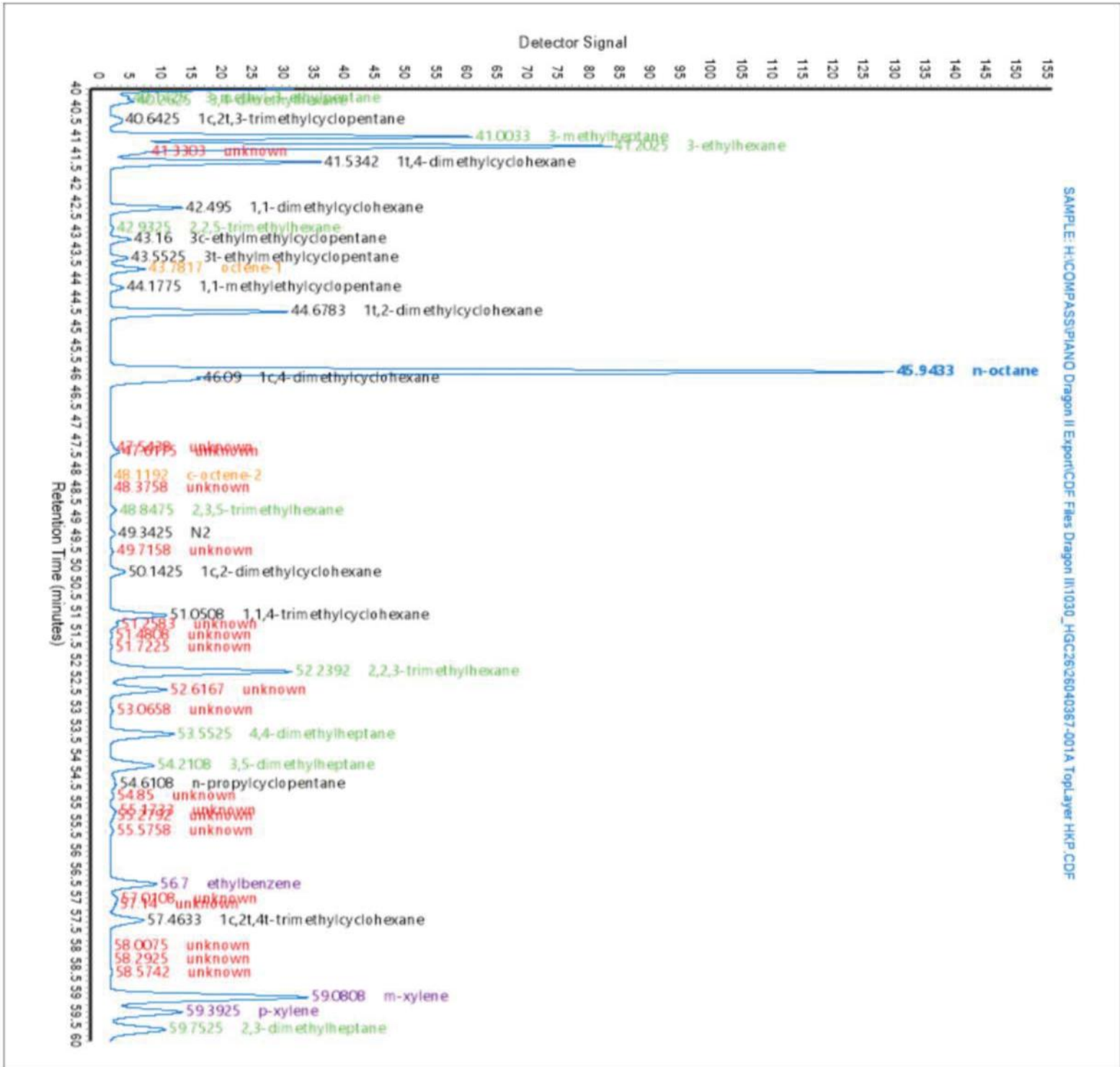
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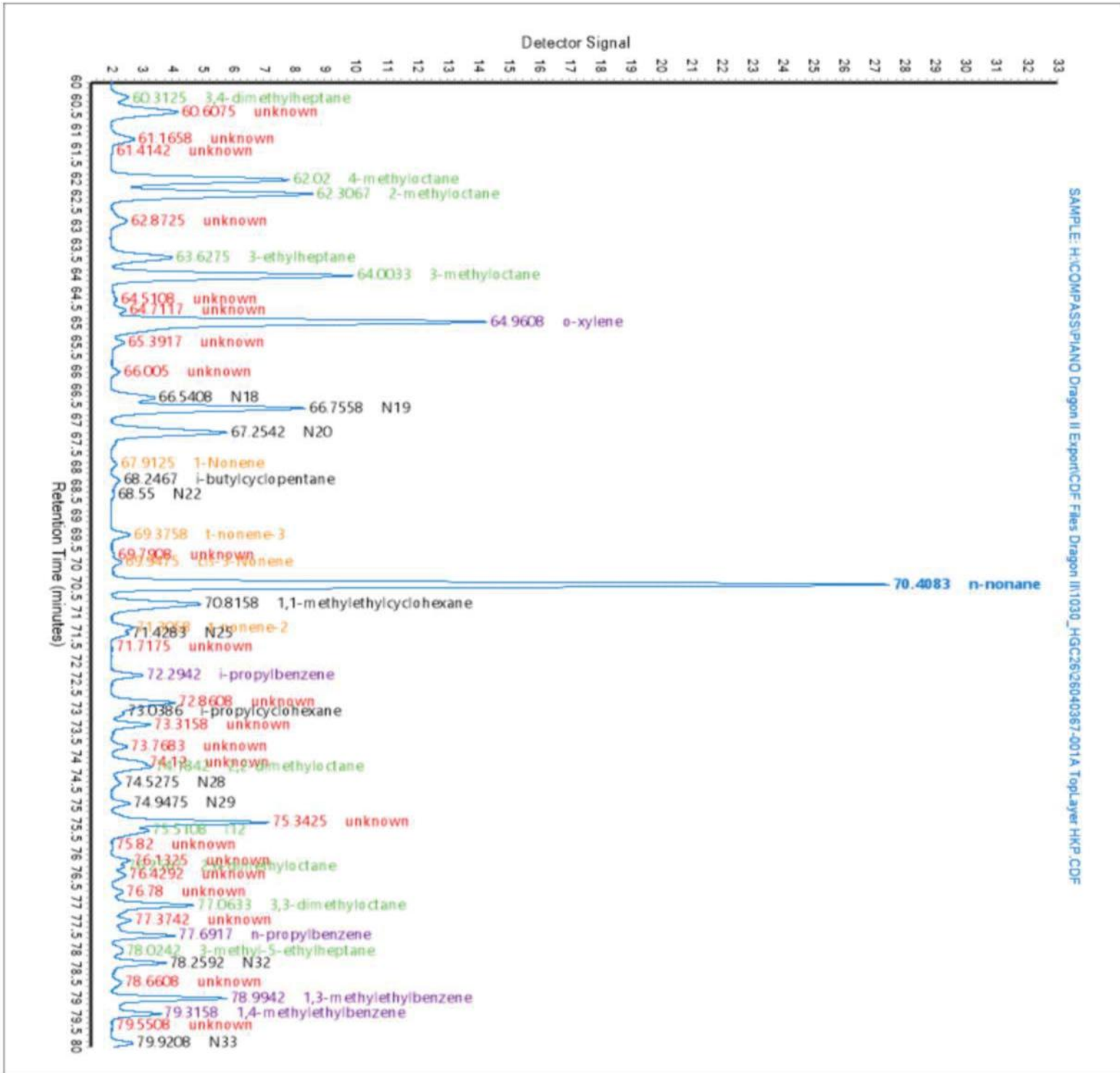
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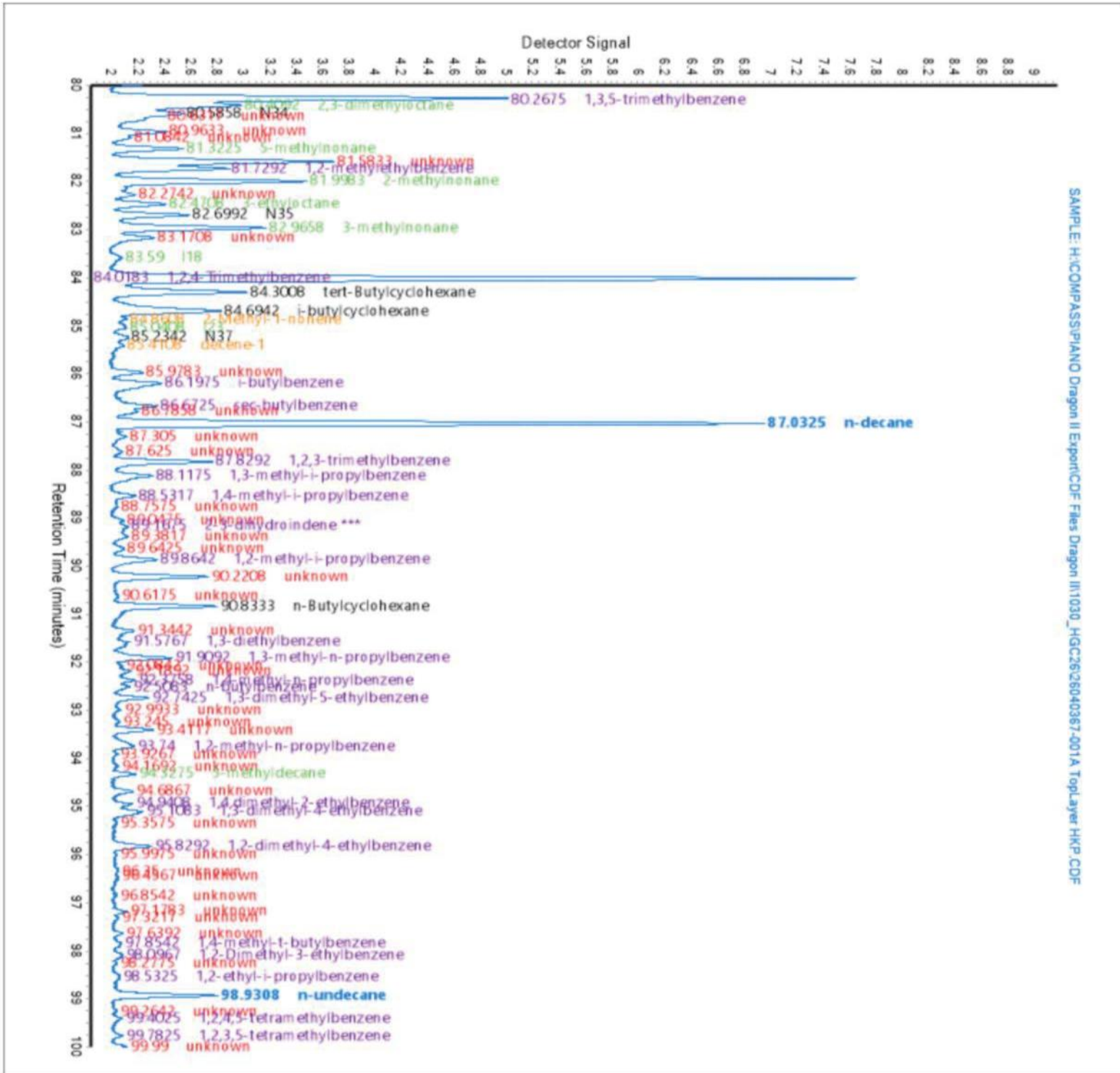
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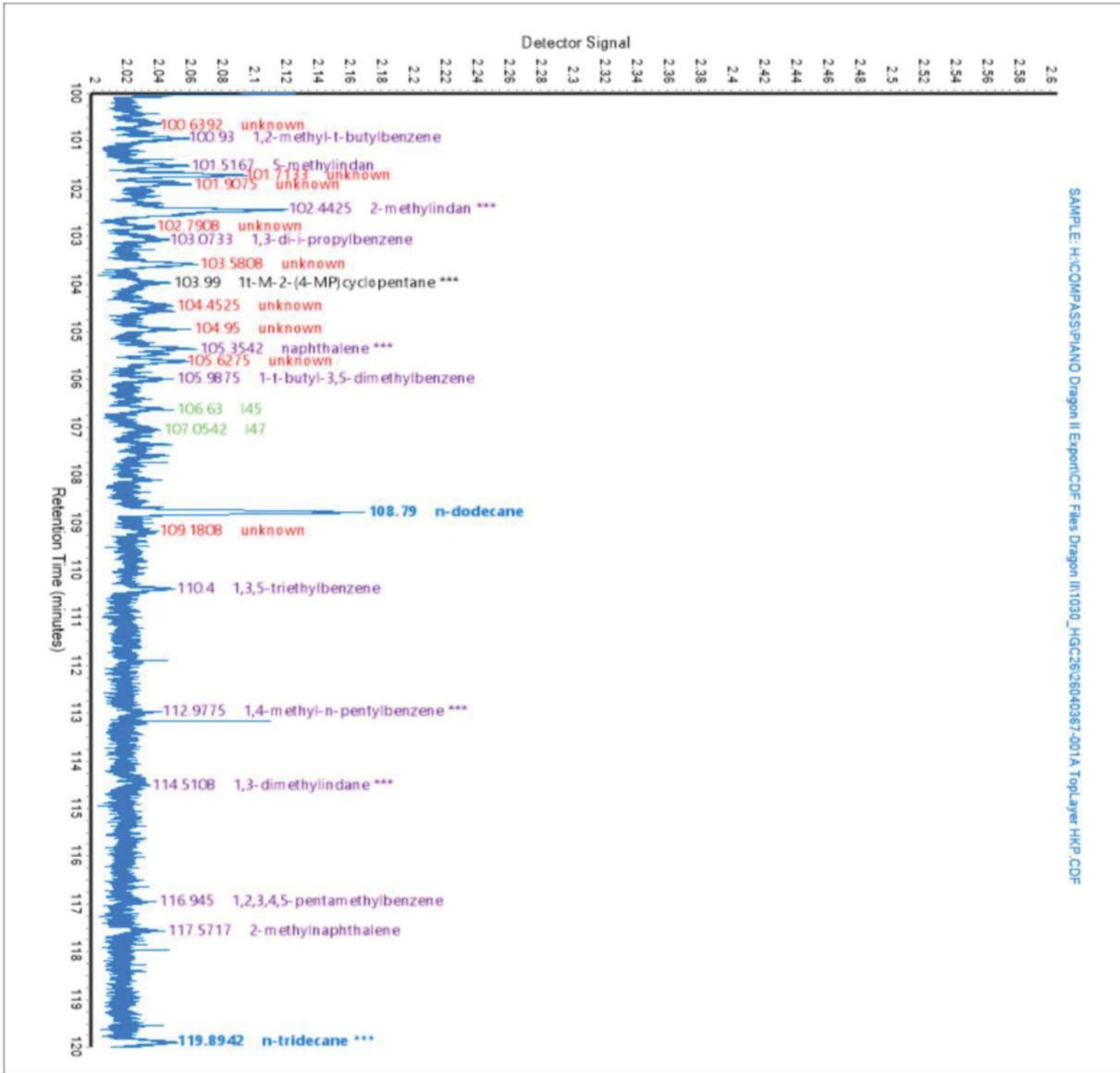
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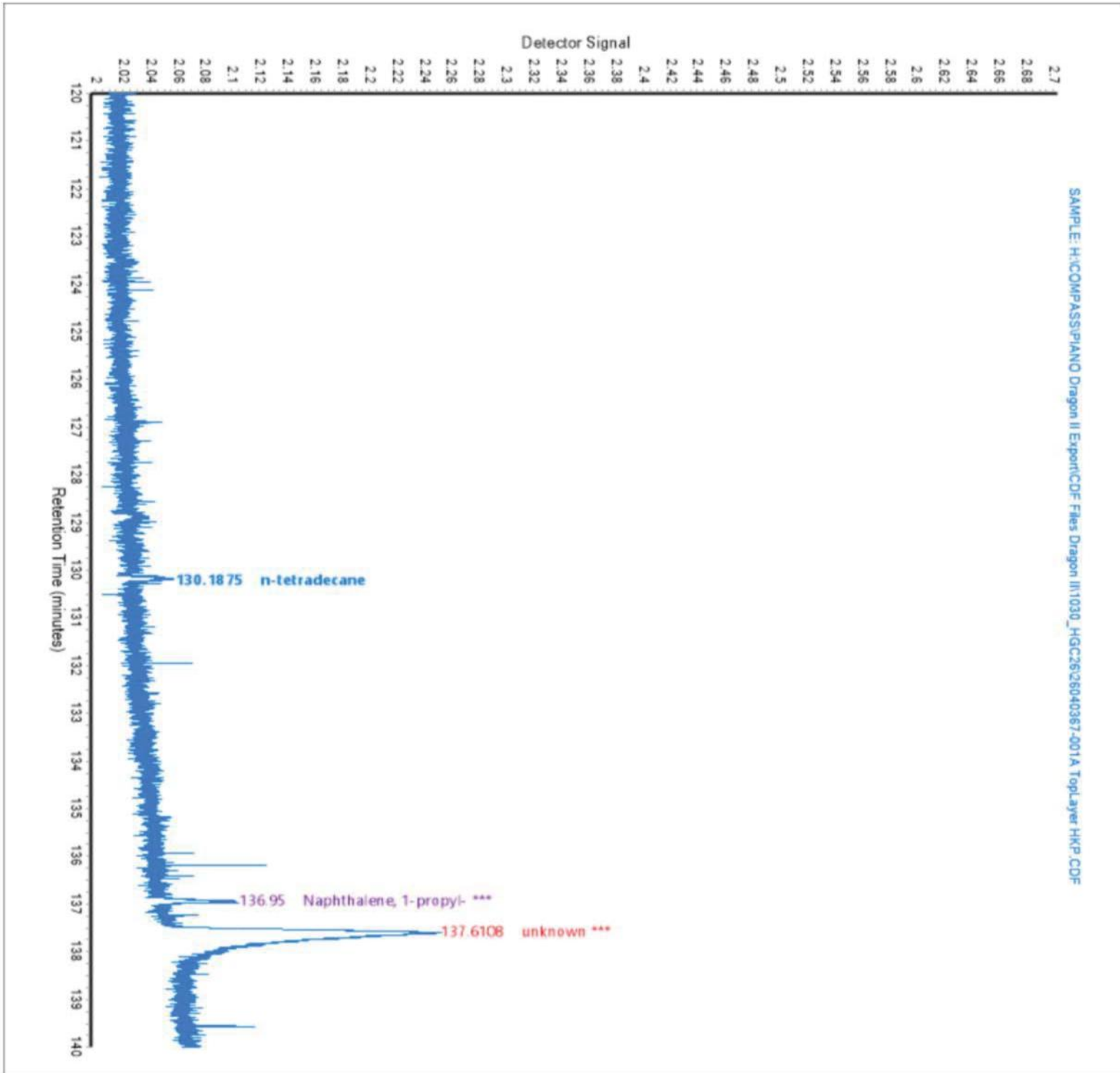
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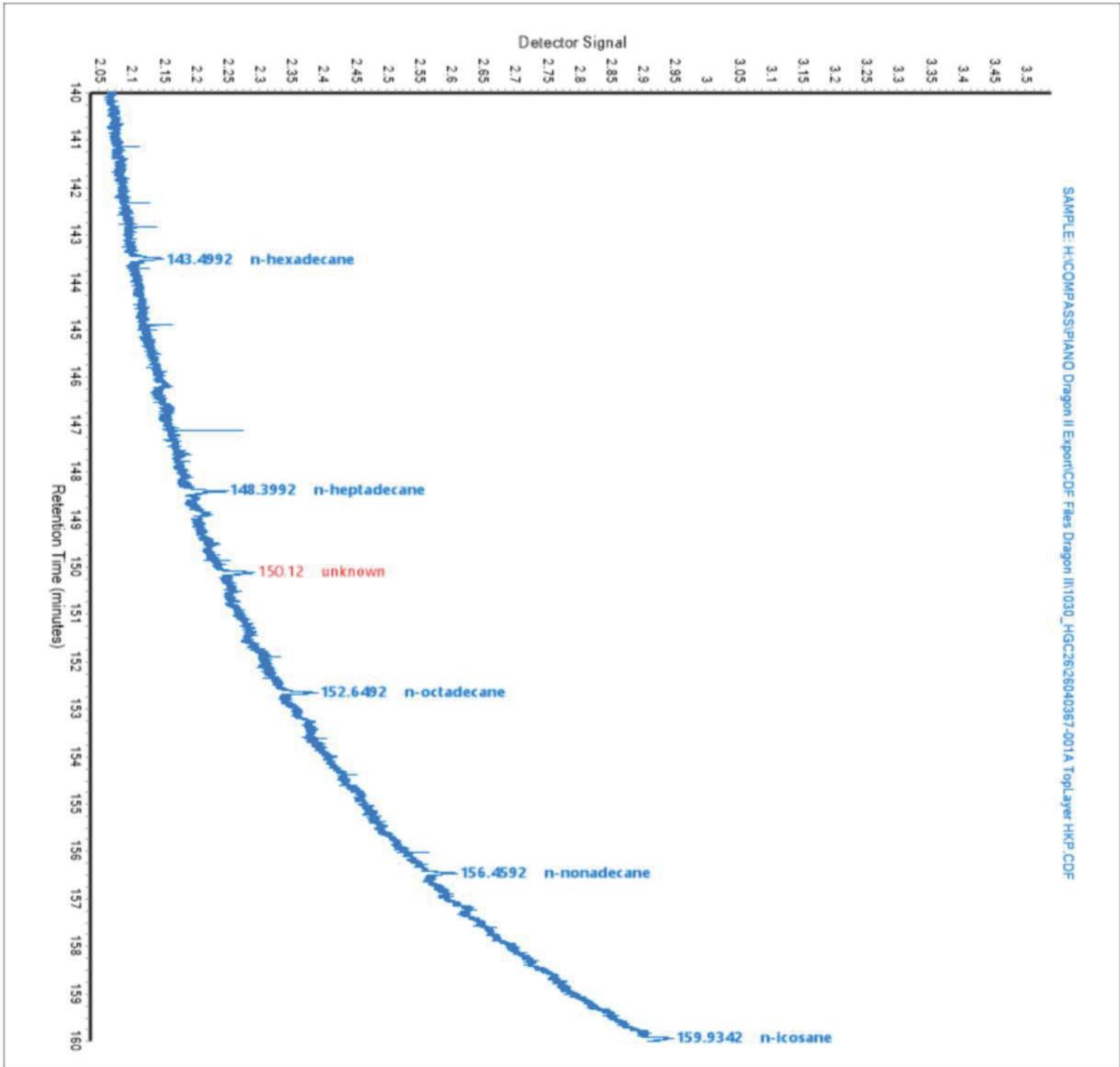
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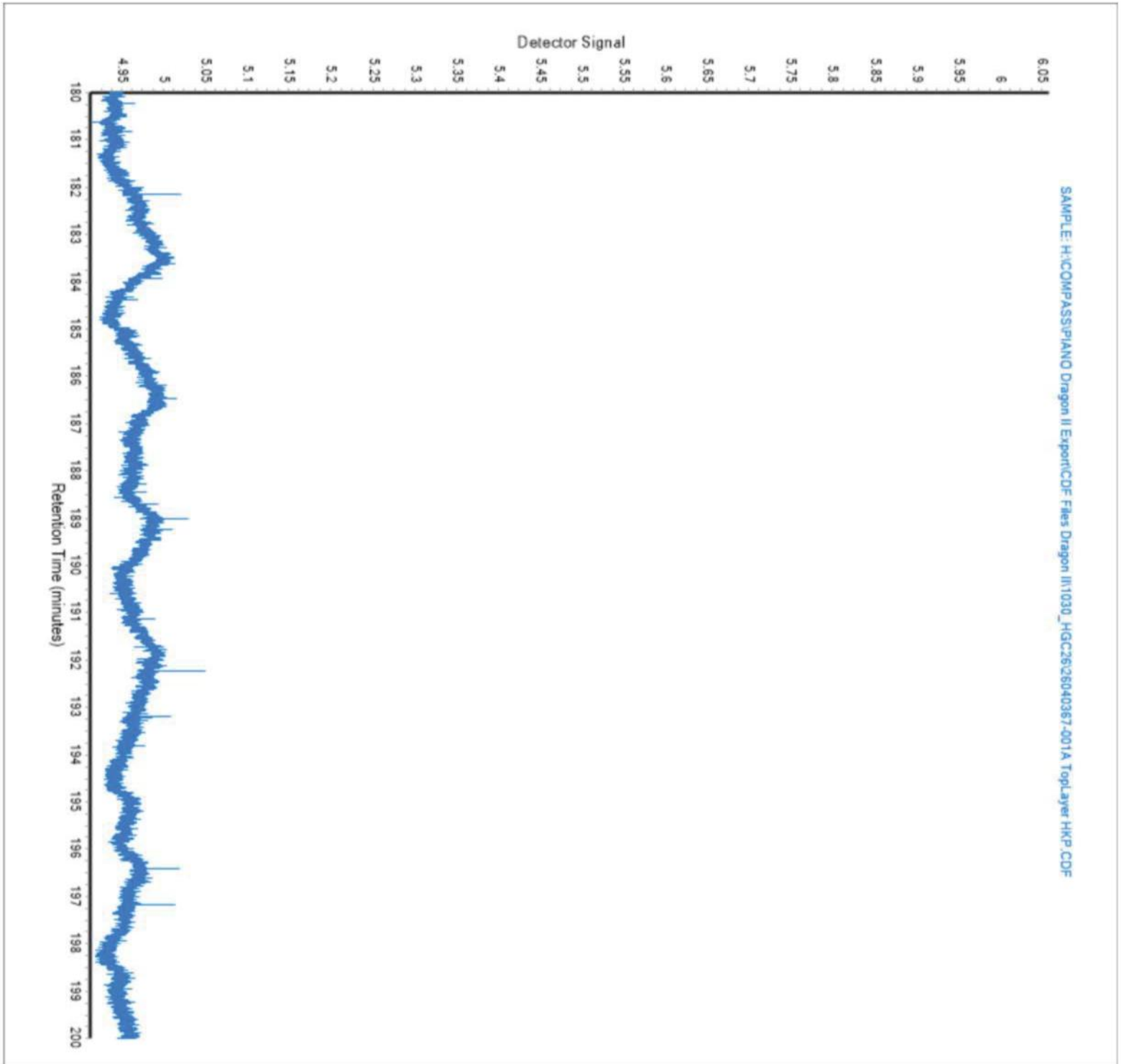
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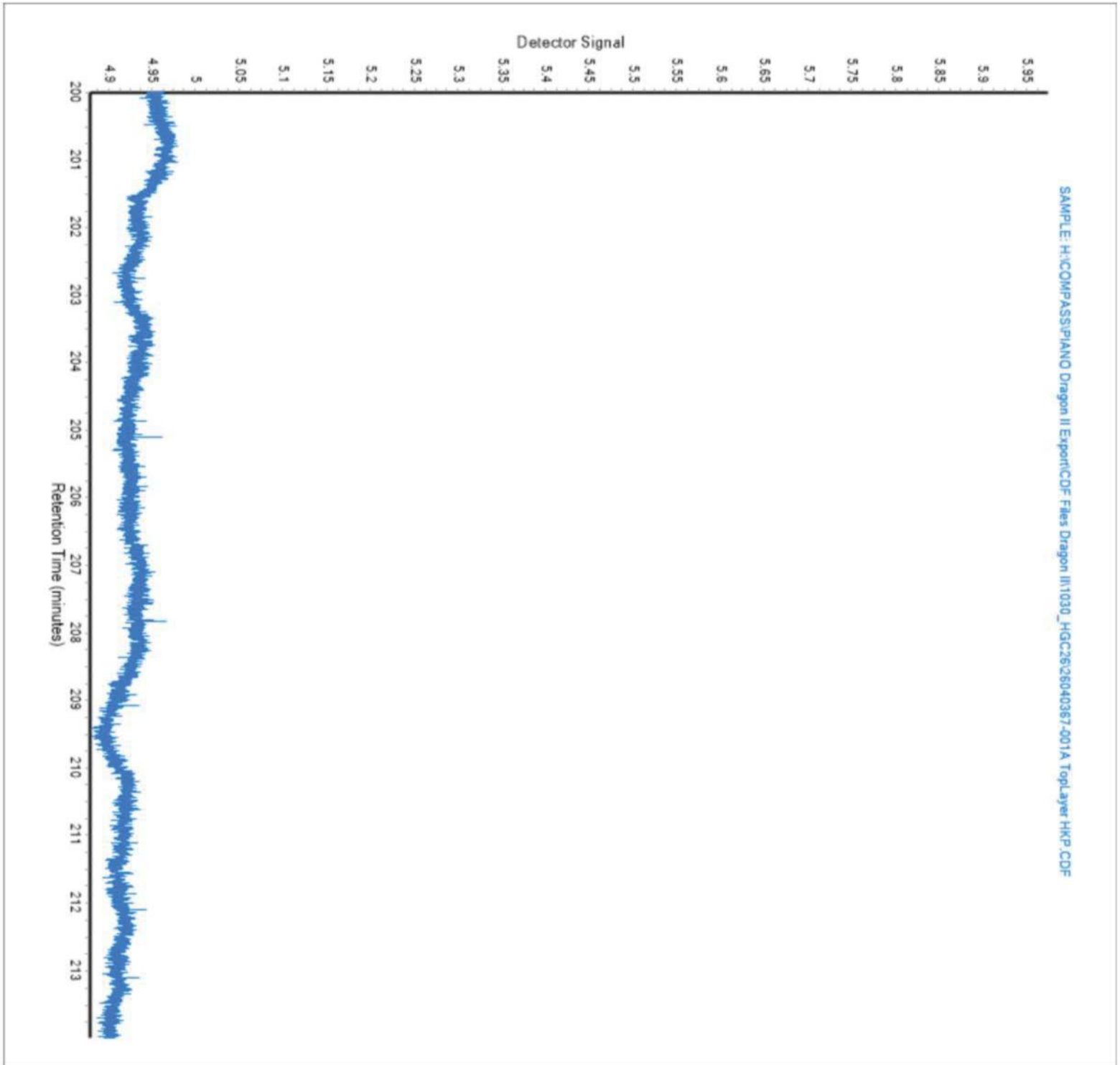
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Certificate of Analysis
 Number: 1030-26040367-001B

Houston Laboratories
 8820 Interchange Drive
 Houston, TX 77054
 Phone 713-660-0901

Luke Arnsberger
 Dolan Intergration Group
 11025 Dover St, Ste 800
 Westminster, CO 80021

Sample ID: 26041347-001A
 Station Name: DCP Wattenburg Pipeline
 Station Number: Spill ID 493177
 Station Location: Hwy 52 & CR19 (G) 2/2026
 Method: Proprietary
 Analyzed: 04/16/2026 by HKP

Report Date: 04/22/2026
 Sampled By:
 Sample Of: Liquid Spot
 Sample Date: 04/01/2026 13:17
 Sample Conditions:
 Received Date: 04/10/2026
 Login Date: 04/13/2026

Analytical Data

Analyte	Result	Units	Detection Limit
Methylamine	ND	wt%	0.01
Dimethylamine	ND	wt%	0.01
Trimethylamine	ND	wt%	0.01
Ethanolamine	14.550	wt%	0.01
2-Dimethylaminoethanol	ND	wt%	0.01
Pyridine	ND	wt%	0.01
Piperazine	ND	wt%	0.01
Diglycolamine	ND	wt%	0.01
DEA + MDEA	ND	wt%	0.01
Butyldiethanolamine	ND	wt%	0.01
Triethanolamine	ND	wt%	0.01

Comments: ND = Not Detected, <0.01 wt%
 Analysis performed on aqueous layer.

Joseph Ponminissery, Laboratory Director

Quality Assurance: The above analyses are performed in accordance with ASTM, UOP, GPA guidelines for quality assurance, unless otherwise stated. The test results apply to the sample as received.



Certificate of Analysis
Number: 1030-26040367-001B

Houston Laboratories
8820 Interchange Drive
Houston, TX 77054
Phone 713-660-0901

Luke Arnsberger
Dolan Intergration Group
11025 Dover St, Ste 800
Westminister, CO 80021

Sample ID: 26041347-001A
Station Name: DCP Wattenburg Pipeline
Station Number: Spill ID 493177
Station Location: Hwy 52 & CR19 (G) 2/2026
Method: Proprietary
Analyzed: 04/16/2026 by HKP

Report Date: 04/22/2026
Sampled By:
Sample Of: Liquid Spot
Sample Date: 04/01/2026 13:17
Sample Conditions:
Received Date: 04/10/2026
Login Date: 04/13/2026

Analytical Data

Analyte	Result	Units	Detection Limit
Ethylene Glycol	0.63	wt%	0.01
Propylene Glycol	ND	wt%	0.01
Diethylene Glycol	ND	wt%	0.01
Triethylene Glycol	70.10	wt%	0.01

Comments: ND = Not Detected, <0.01 wt%
Analysis performed on aqueous layer.

Joseph Ponminissery, Laboratory Director

Quality Assurance: The above analyses are performed in accordance with ASTM, UOP, GPA guidelines for quality assurance, unless otherwise stated. The test results apply to the sample as received.



Certificate of Analysis
 Number: 1030-26040367-001B

Houston Laboratories
 8820 Interchange Drive
 Houston, TX 77054
 Phone 713-660-0901

Luke Arnsberger
 Dolan Intergration Group
 11025 Dover St, Ste 800
 Westminster, CO 80021

Sample ID: 26041347-001A
 Station Name: DCP Wattenburg Pipeline
 Station Number: Spill ID 493177
 Station Location: Hwy 52 & CR19 (G) 2/2026
 Method: ORYX
 Analyzed: 04/21/2026 by HKP

Report Date: 04/22/2026
 Sampled By:
 Sample Of: Liquid Spot
 Sample Date: 04/01/2026 13:17
 Sample Conditions:
 Received Date: 04/10/2026
 Login Date: 04/13/2026

Analytical Data

Analyte	Result	Units	Detection Limit
Methane	ND	wt%	0.001
Ethane	ND	wt%	0.001
Propane	ND	wt%	0.001
Iso-Butane	ND	wt%	0.001
n-Butane	0.969	wt%	0.001
Iso-Pentane	2.656	wt%	0.001
n-Pentane	5.251	wt%	0.001
Cyclopentane	0.984	wt%	0.001
i-Hexanes	8.112	wt%	0.001
n-Hexane	9.633	wt%	0.001
Methylcyclopentane	4.763	wt%	0.001
Benzene	4.994	wt%	0.001
Cyclohexane	4.394	wt%	0.001
i-Heptanes	9.614	wt%	0.001
2,2,4-Trimethylpentane	ND	wt%	0.001
n-Heptane	7.230	wt%	0.001
Methylcyclohexane	7.435	wt%	0.001
Toluene	6.637	wt%	0.001
i-Octanes	9.600	wt%	0.001
n-Octane	2.664	wt%	0.001
Ethylbenzene	0.274	wt%	0.001
m-Xylene	1.426	wt%	0.001
p-Xylene	0.466	wt%	0.001
o-Xylene	0.530	wt%	0.001
i-Nonanes	3.446	wt%	0.001
n-Nonane	0.582	wt%	0.001
i-Decanes	1.072	wt%	0.001
n-Decane	0.085	wt%	0.001
Undecanes Plus	7.182	wt%	0.001

Comments: Undecanes Plus hydrocarbon range is C11-C26.
 Major components in Undecanes Plus are C15, C22.
 Analysis performed on aqueous layer.

Joseph Ponminissery, Laboratory Director

Quality Assurance: The above analyses are performed in accordance with ASTM, UOP, GPA guidelines for quality assurance, unless otherwise stated. The test results apply to the sample as received.