LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA N\DATA\BN032318\
Data File : BN000577.D
Acq On : 23 Mar 2018 14:46
Operator : JU/SJ
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.349	489	495	500	rBV2	41492	62793	1.08%	0.242%
2	6.855	573	581	586	rBV2	23676	42987	0.74%	0.165%
3	7.002	602	606	614	rVB3	13172	33541	0.58%	0.129%
4	7.355	662	666	671	rVB2	28074	41574	0.72%	0.160%
5	7.413	671	676	680	rBV	29094	42983	0.74%	0.165%
6	7.466	680	685	689	rBV	29539	41560	0.72%	0.160%
7	7.525	689	695	701	rVB2	53125	82682	1.42%	0.318%
8	7.607	701	709	712	rBV	18644	30306	0.52%	0.117%
9	7.802	734	742	751	rVB2	49719	106003	1.82%	0.408%
10	8.002	770	776	780	rBV	144516	210856	3.63%	0.811%
11	8.043	780	783	788	rVB	36276	47429	0.82%	0.183%
12	8.360	831	837	840	rBV	31220	43483	0.75%	0.167%
13	8.413	840	846	850	rBV	628753	1055715	18.17%	4.062%
14	8.454	850	853	860	rVB	223576	315916	5.44%	1.216%
15	8.519	860	864	870	rBV4	25564	46277	0.80%	0.178%
16	8.602	870	878	883	rBV	105519	180517	3.11%	0.695%
17	8.837	912	918	926	rBV	98517	183209	3.15%	0.705%
18	8.919	926	932	935	rVV3	25262	50125	0.86%	0.193%
19	8.978	935	942	949	rVB3	30449	69565	1.20%	0.268%
20	9.049	949	954	961	rBV2	68240	106908	1.84%	0.411%
21	9.160	967	973	977	rBV	35260	58737	1.01%	0.226%
22	9.213	977	982	986	rVB2	16495	29096	0.50%	0.112%
23	9.366	1003	1008	1011	rBV	17888	28031	0.48%	0.108%
24	9.419	1014	1017	1022	rVB	17022	27188	0.47%	0.105%
25	9.519	1029	1034	1042	rVB3	35746	63907	1.10%	0.246%
26	9.625	1044	1052	1057	rBV	185543	283484	4.88%	1.091%
27	9.678	1057	1061	1067	rBV	30676	50978	0.88%	0.196%
28	9.849	1084	1090	1097	rBV6	16616	52632	0.91%	0.203%
29	10.201	1146	1150	1155	rVB	23448	38201	0.66%	0.147%
30	10.278	1155	1163	1164	rBV3	16579	35108	0.60%	0.135%
31	10.349	1170	1175	1182	rVB4	36695	62251	1.07%	0.240%
32	10.478	1190	1197	1202	rVB2	26435	59214	1.02%	0.228%
33	10.549	1202	1209	1218	rBV6	19578	44841	0.77%	0.173%
34	10.631	1218	1223	1228	rVV3	29313	46608	0.80%	0.179%

LSC Area Percent Report

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Acq On : 23 Mar 2018 14:46
Operator : JU/SJ
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Title : SVOA CALIBRATION

35	10.731	1236	1240	1246	rVV4	18297	33446	0.58%	0.129%
36	11.196	1307	1319	1323	rVV3	113168	343064	5.91%	1.320%
37	11.248	1323	1328	1337	rVV	884699	1567193	26.98%	6.031%
38	11.325	1337	1341	1344	rVV3	47497	80816	1.39%	0.311%
39	11.360	1344	1347	1353	rVB3	32875	57082	0.98%	0.220%
40	11.590	1380	1386	1394	rVB	105518	168536	2.90%	0.649%
41	11.719	1400	1408	1414	rBV	32911	54818	0.94%	0.211%
42	11.925	1435	1443	1456	rVV10	19578	65553	1.13%	0.252%
43	12.107	1468	1474	1480	rVV4	26765	58384	1.00%	0.225%
44	12.219	1487	1493	1498	rVV3	37300	73785	1.27%	0.284%
45	12.307	1504	1508	1516	rVB	50561	78351	1.35%	0.302%
46	12.454	1528	1533	1543	rVB2	81680	128653	2.21%	0.495%
47	12.601	1553	1558	1562	rBV	83407	130130	2.24%	0.501%
48	12.648	1562	1566	1572	rVB	62821	89991	1.55%	0.346%
49	12.737	1572	1581	1588	rBV5	22636	70186	1.21%	0.270%
50	13.013	1623	1628	1631	rBV4	25231	43365	0.75%	0.167%
51	13.090	1636	1641	1651	rVB3	10639	30477	0.52%	0.117%
52	13.237	1658	1666	1669	rBV6	17073	41756	0.72%	0.161%
53	13.395	1686	1693	1696	rBV7	13994	29441	0.51%	0.113%
54	13.537	1713	1717	1723	rBV2	32386	48156	0.83%	0.185%
55	13.754	1749	1754	1760	rVB2	32380	47373	0.82%	0.182%
56	13.819	1760	1765	1776	rBV	103256	172637	2.97%	0.664%
57	13.966	1787	1790	1800	rVB6	11567	30609	0.53%	0.118%
58	14.066	1800	1807	1810	rBV4	26221	44845	0.77%	0.173%
59	14.348	1849	1855	1863	rBV10	28768	74738	1.29%	0.288%
60	14.472	1868	1876	1880	rVB5	41762	88385	1.52%	0.340%
61	14.637	1898	1904	1913	rVB6	39021	98372	1.69%	0.379%
62	14.731	1915	1920	1925	rBV3	19930	38920	0.67%	0.150%
63	14.878	1941	1945	1951	rVB	342734	446627	7.69%	1.719%
64	15.031	1964	1971	1984	rVB2	1339608	2097970	36.11%	8.073%
65	15.142	1984	1990	1995	rBV	105055	141523	2.44%	0.545%
66	15.307	2013	2018	2025	rBV4	28839	60863	1.05%	0.234%
67	15.401	2031	2034	2044	rVB6	28959	64761	1.11%	0.249%
68	15.489	2044	2049	2053	rBV3	34925	53236	0.92%	0.205%
69	15.642	2067	2075	2087	rBV	360978	642661	11.06%	2.473%
70	15.760	2090	2095	2099	rVV	107798	150984	2.60%	0.581%
71	15.819	2099	2105	2111	rVB	126086	221304	3.81%	0.852%

LSC Area Percent Report

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Acq On : 23 Mar 2018 14:46
Operator : JU/SJ
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Title : SVOA CALIBRATION

72	16.225	2165	2174	2177	rBV	44694	85295	1.47%	0.328%
73	16.266	2177	2181	2186	rVB4	20209	31362	0.54%	0.121%
74	16.607	2235	2239	2246	rVB4	43244	67937	1.17%	0.261%
75	16.678	2246	2251	2254	rBV	131662	203012	3.49%	0.781%
76	16.813	2271	2274	2280	rBV6	15457	30707	0.53%	0.118%
77	16.919	2285	2292	2297	rBV9	15633	33785	0.58%	0.130%
78	17.042	2310	2313	2316	rVV3	27268	37244	0.64%	0.143%
79	17.407	2368	2375	2381	rBV	3970402	5809636	100.00%	22.356%
80	17.460	2381	2384	2389	rVV	163067	251608	4.33%	0.968%
81	17.513	2389	2393	2409	rVB3	65596	162303	2.79%	0.625%
82	17.754	2428	2434	2442	rBV2	1643794	2456315	42.28%	9.452%
83	18.001	2471	2476	2478	rBV4	21905	33505	0.58%	0.129%
84	18.195	2504	2509	2516	rVB	103414	132750	2.28%	0.511%
85	18.366	2535	2538	2544	rVB2	20860	29744	0.51%	0.114%
86	18.872	2620	2624	2629	rVB	74919	89437	1.54%	0.344%
87	19.495	2726	2730	2732	rBV	59793	76046	1.31%	0.293%
88	19.519	2732	2734	2737	rVB2	33088	33263	0.57%	0.128%
89	20.060	2822	2826	2831	rBV	42322	48330	0.83%	0.186%
90	20.589	2912	2916	2920	rBV2	39528	47582	0.82%	0.183%
91	21.077	2996	2999	3003	rBV2	36531	39313	0.68%	0.151%
92	21.542	3074	3078	3082	rBV	57630	66855	1.15%	0.257%
93	21.883	3131	3136	3145	rBV2	1789248	2470731	42.53%	9.508%
94	21.989	3151	3154	3161	rVB3	35785	45263	0.78%	0.174%
95	22.460	3230	3234	3242	rVB	51288	74703	1.29%	0.287%
96	22.965	3317	3320	3324	rBV2	42169	57227	0.99%	0.220%
97	23.530	3412	3416	3423	rVB	56242	88685	1.53%	0.341%
98	24.160	3519	3523	3531	rVB	60045	110136	1.90%	0.424%
99	24.371	3551	3559	3572	rVB2	933141	1934122	33.29%	7.443%
100	24.889	3643	3647	3652	rVB2	53505	96307	1.66%	0.371%

Sum of corrected areas: 25986899

Data Path : Z:\HPCHEM1\BNA N\DATA\BN032318\
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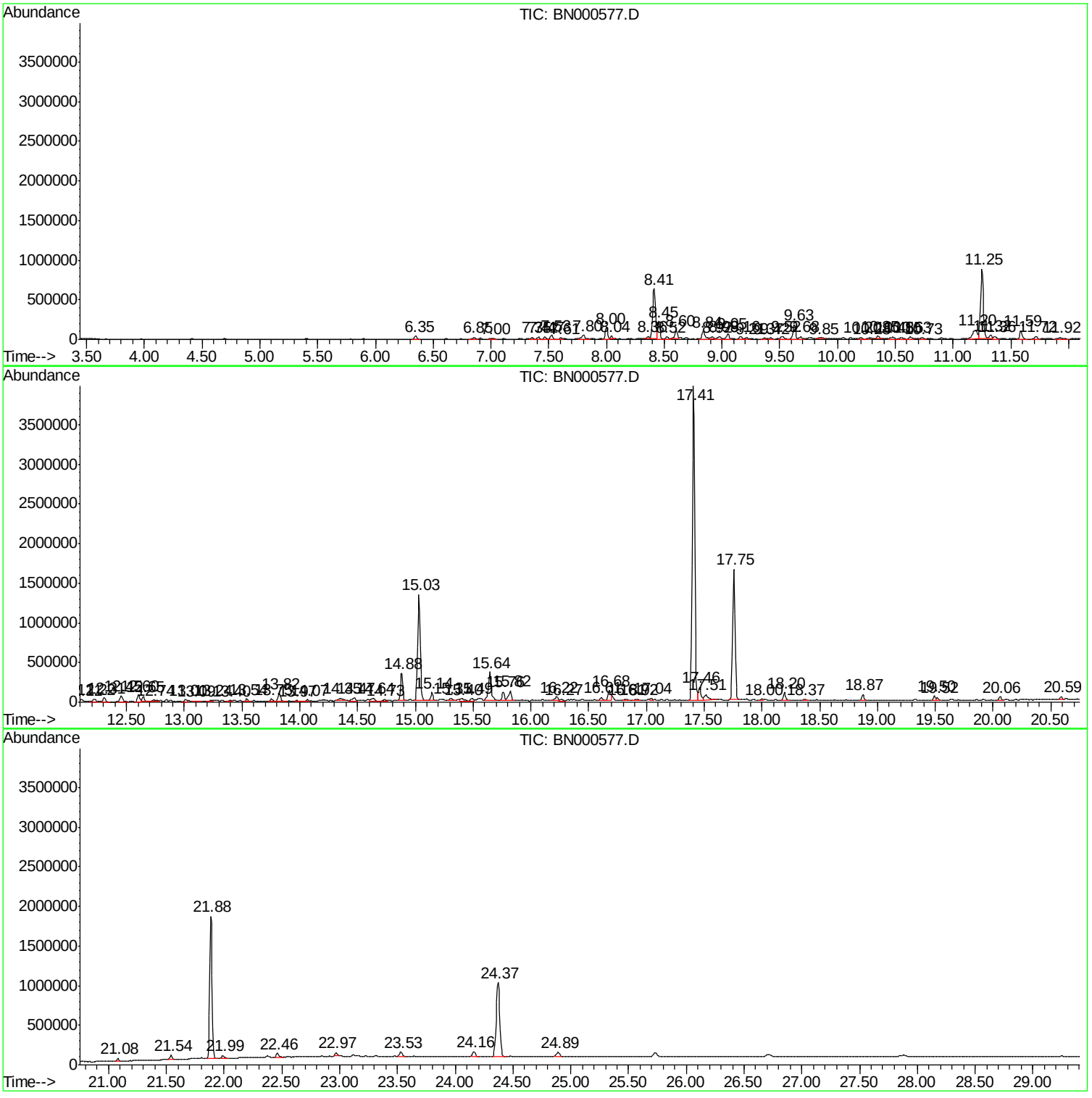
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA N\DATA\BN032318\
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Sample : J1905-07DL3 500X
Misc :
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Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

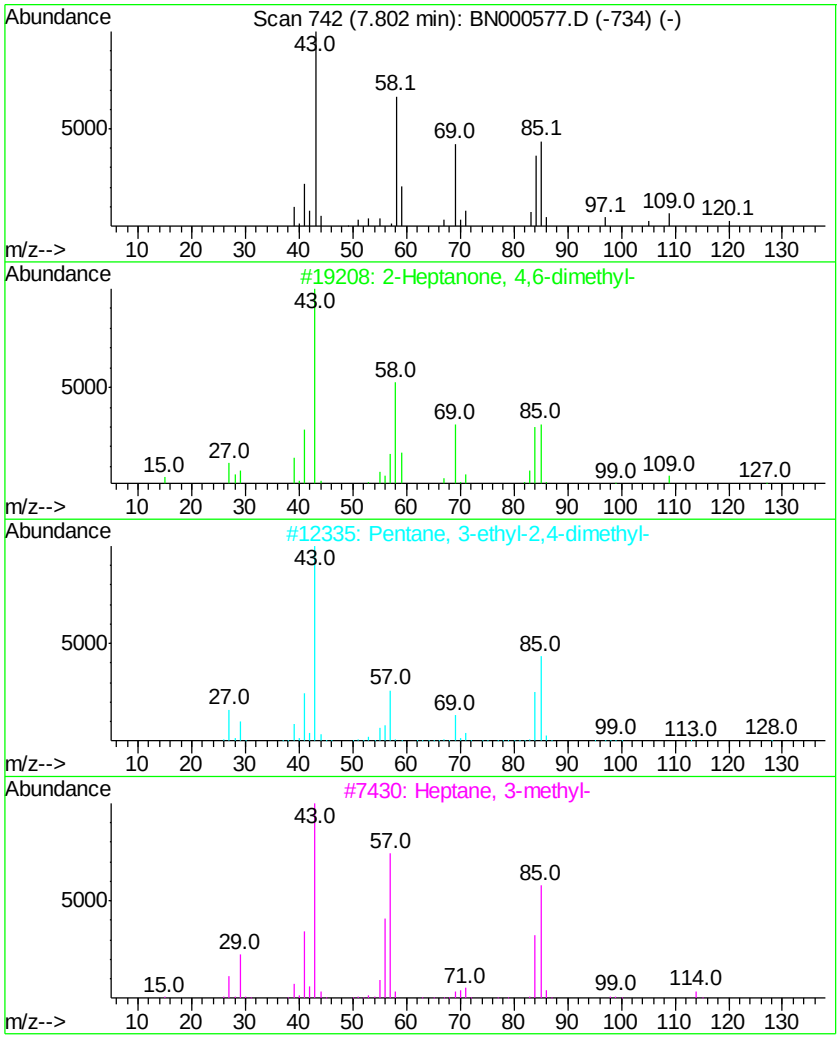
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

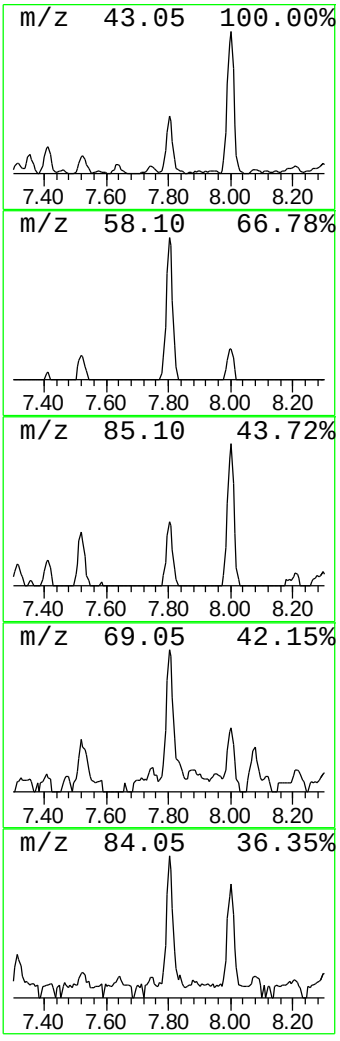
****

Peak Number 1 2-Heptanone, 4,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	2.01 ng/ul	106003	1,4-Dichlorobenzene-d4	8.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	64
2	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	43
3	Heptane, 3-methyl-	114	C8H18	000589-81-1	38
4	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	37
5	Hexane, 3-ethyl-	114	C8H18	000619-99-8	32





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Acq On : 23 Mar 2018 14:46
Operator : JU/SJ
Sample : J1905-07DL3 500X
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ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

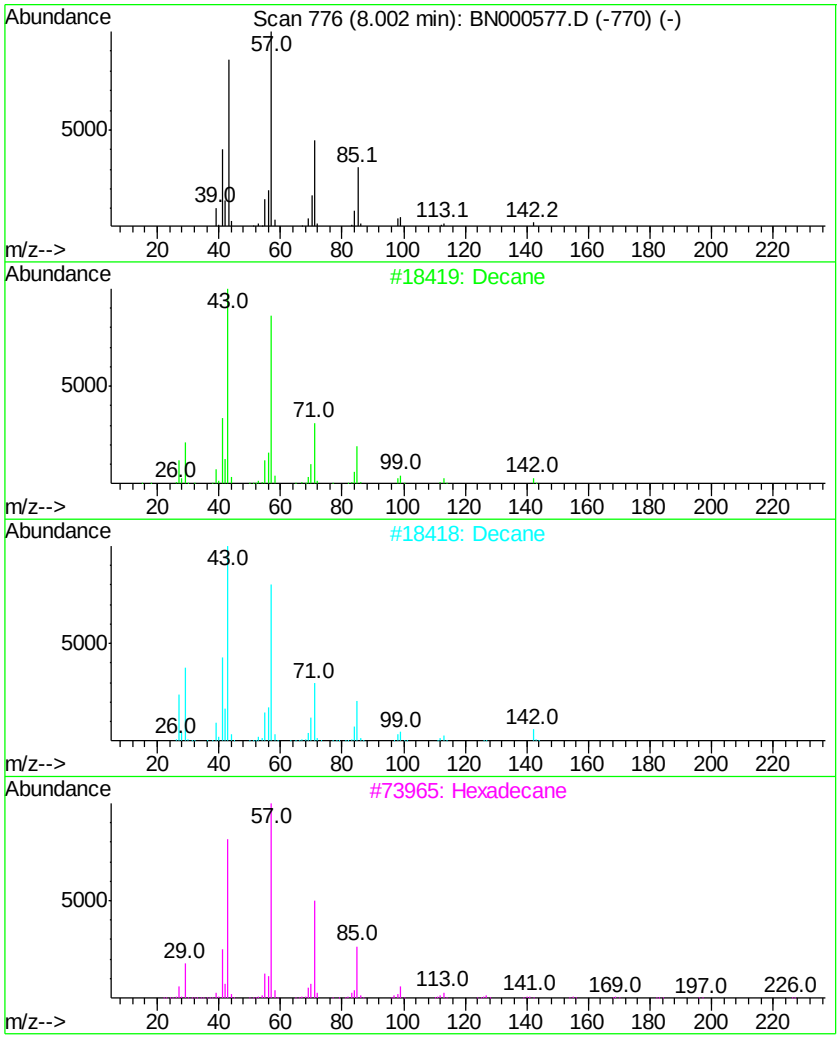
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

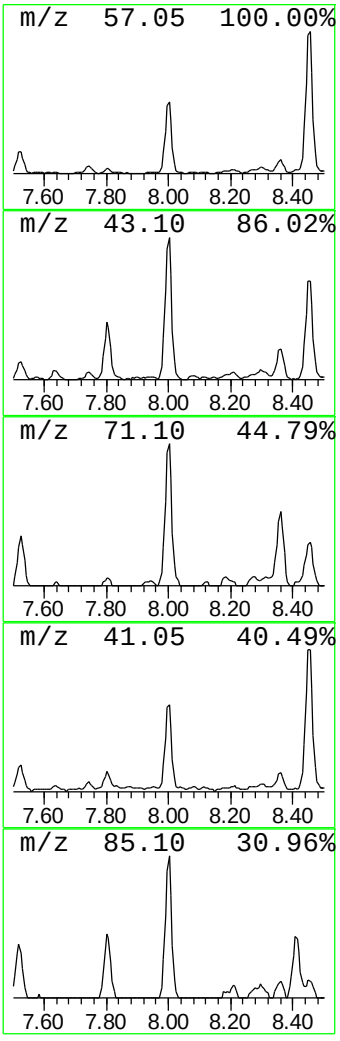
****

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.00	3.99 ng/ul	210856	1,4-Dichlorobenzene-d4	8.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	91
2	Decane	142	C10H22	000124-18-5	90
3	Hexadecane	226	C16H34	000544-76-3	72
4	Nonane	128	C9H20	000111-84-2	72
5	Nonane	128	C9H20	000111-84-2	72





Library Search Compound Report

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Operator : JU/SJ
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Misc :
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Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

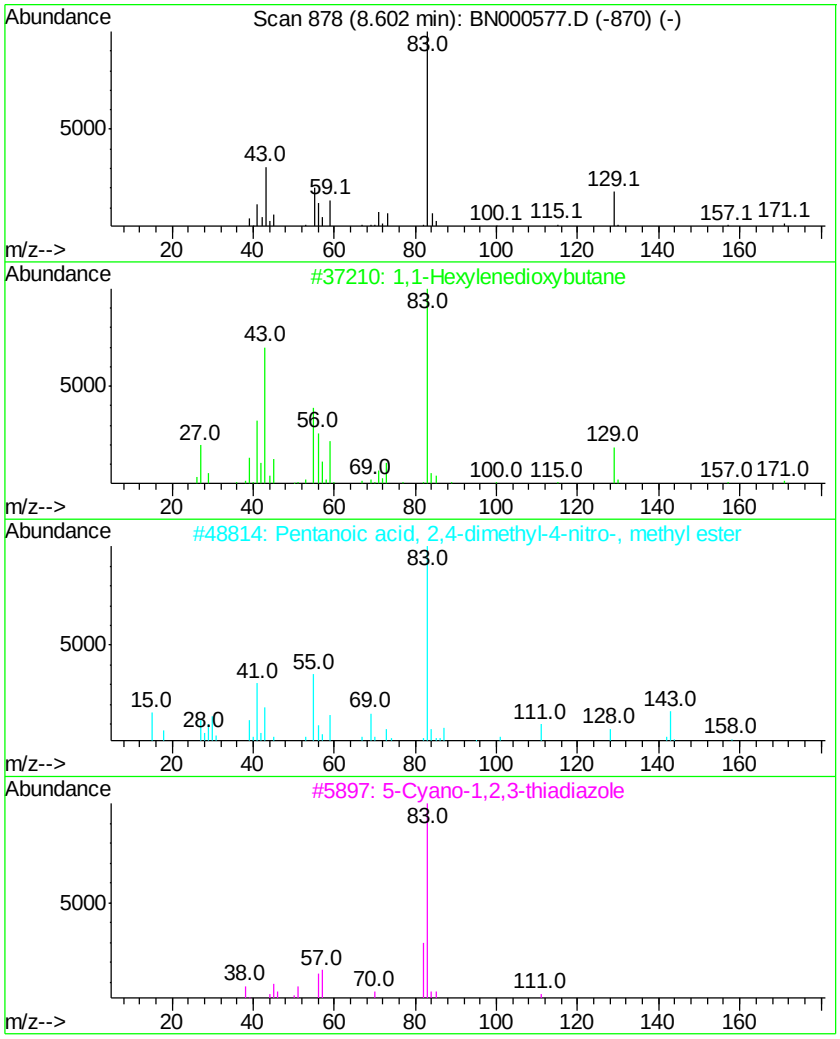
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

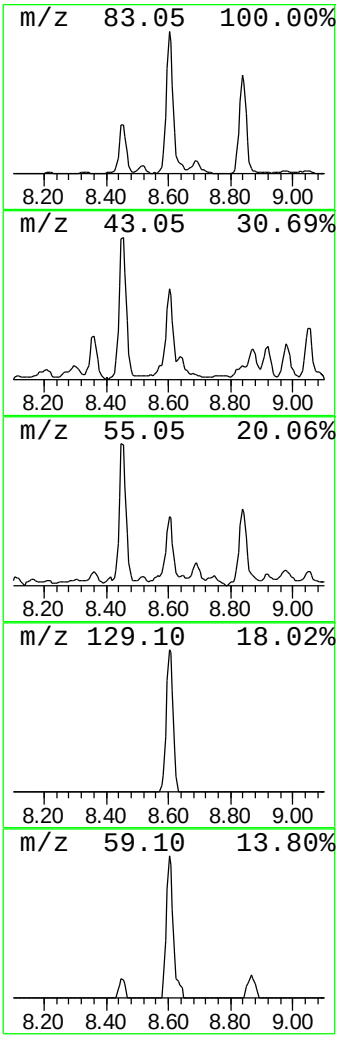
****

Peak Number 3 1,1-Hexylenedioxybutane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.60	3.42 ng/ul	180517	1,4-Dichlorobenzene-d4	8.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	50
2	Pentanoic acid, 2,4-dimethyl-4-n...	189	C8H15NO4	005762-40-3	40
3	5-Cyano-1,2,3-thiadiazole	111	C3HN3S	057352-02-0	38
4	Cyclohexane, undecyl-	238	C17H34	054105-66-7	38
5	3-Methyl-2-butenoic acid, 2,2-di...	170	C10H18O2	1000279-23-0	36





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Acq On : 23 Mar 2018 14:46
Operator : JU/SJ
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

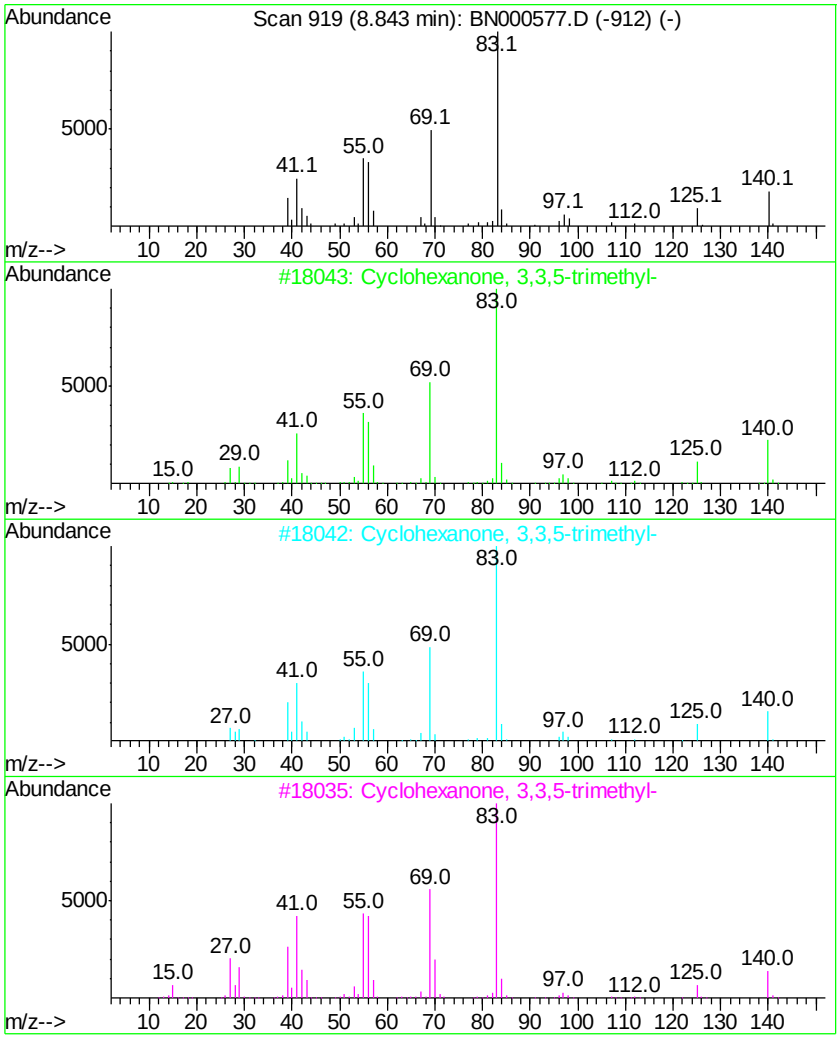
Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

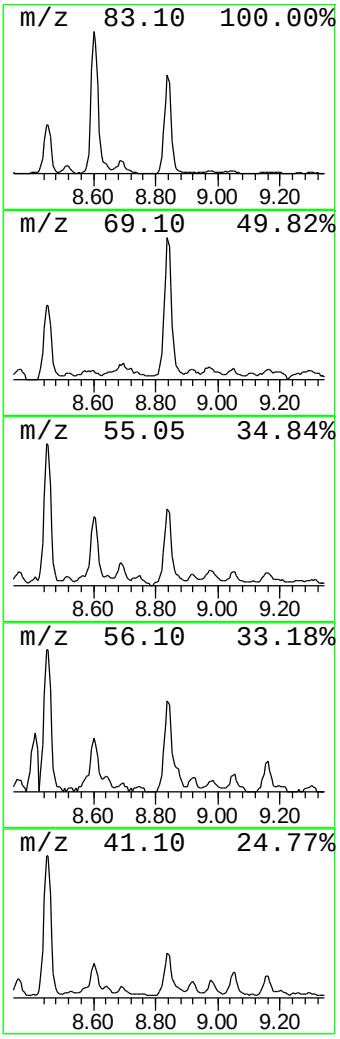
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclohexanone, 3,3,5-trimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.84	3.47 ng/ul	183209	1,4-Dichlorobenzene-d4	8.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	97
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4	Cyclohexane, butyl-	140	C10H20	001678-93-9	72
5	Cyclohexane, butyl-	140	C10H20	001678-93-9	64





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA N\DATA\BN032318\
Data File : BN000577.D
Acq On : 23 Mar 2018 14:46
Operator : JU/SJ
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

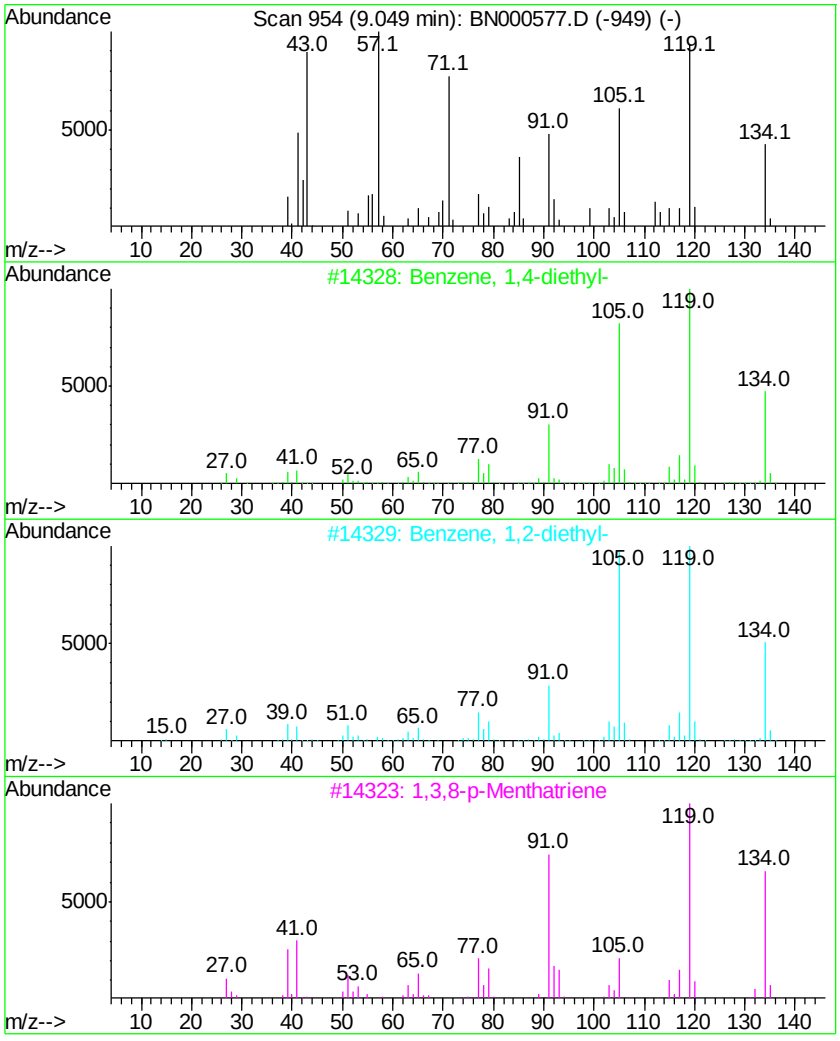
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

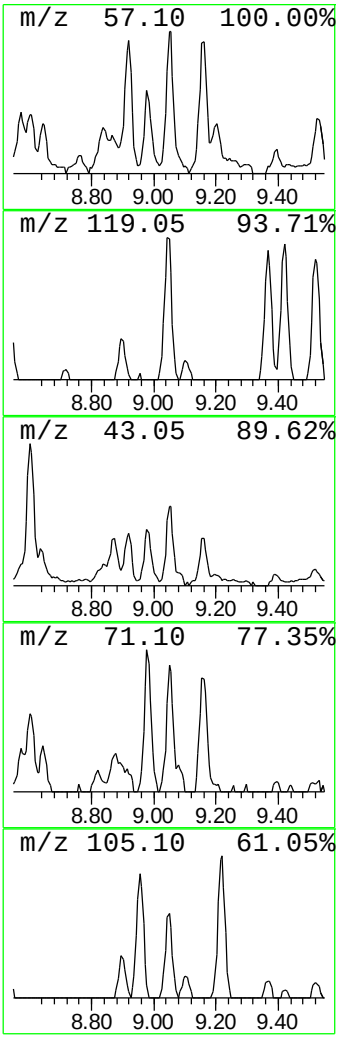
****

Peak Number 5 Benzene, 1,4-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	2.03 ng/ul	106908	1,4-Dichlorobenzene-d4	8.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	89
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	86
3	1,3,8-p-Menthatriene	134	C10H14	021195-59-5	70
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	55
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	55





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032318\
Data File : BN000577.D
Acq On : 23 Mar 2018 14:46
Operator : JU/SJ
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

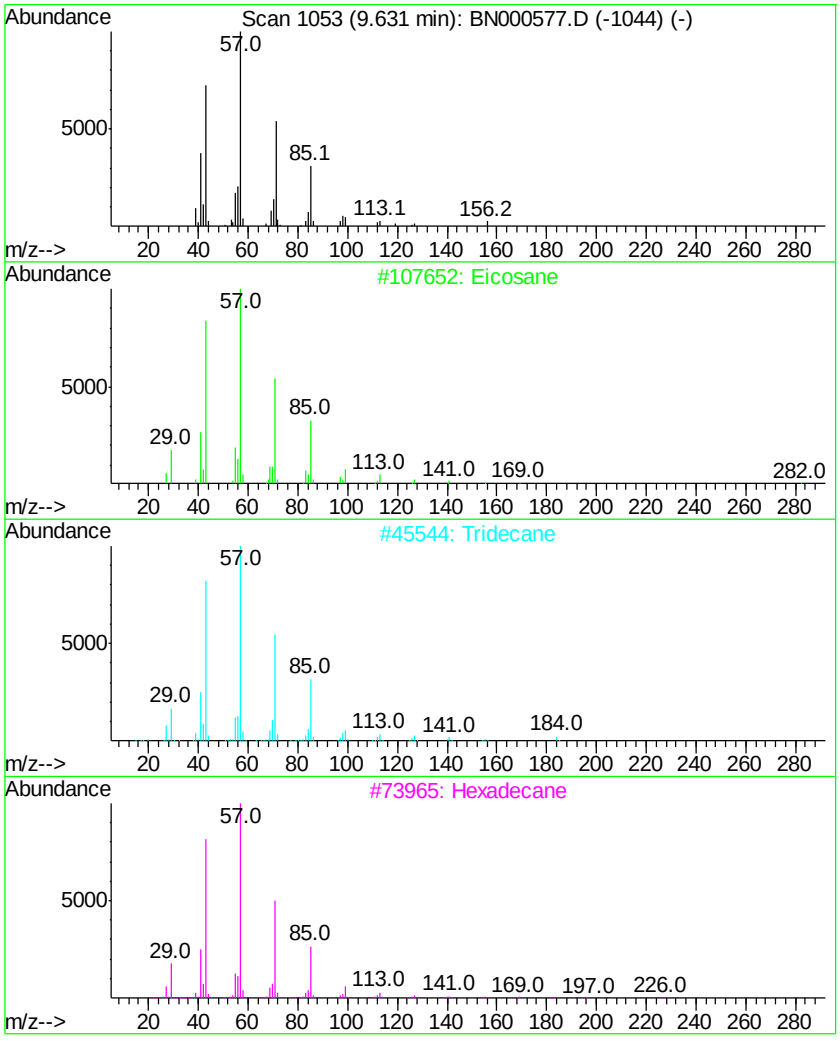
Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

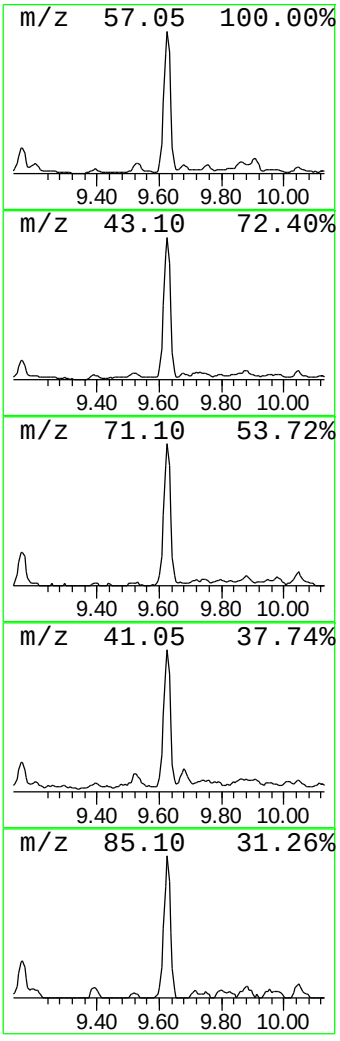
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	5.37 ng/ul	283484	1,4-Dichlorobenzene-d4	8.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	72
2	Tridecane	184	C13H28	000629-50-5	72
3	Hexadecane	226	C16H34	000544-76-3	59
4	Decane, 5-propyl-	184	C13H28	017312-62-8	59
5	Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	59





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Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

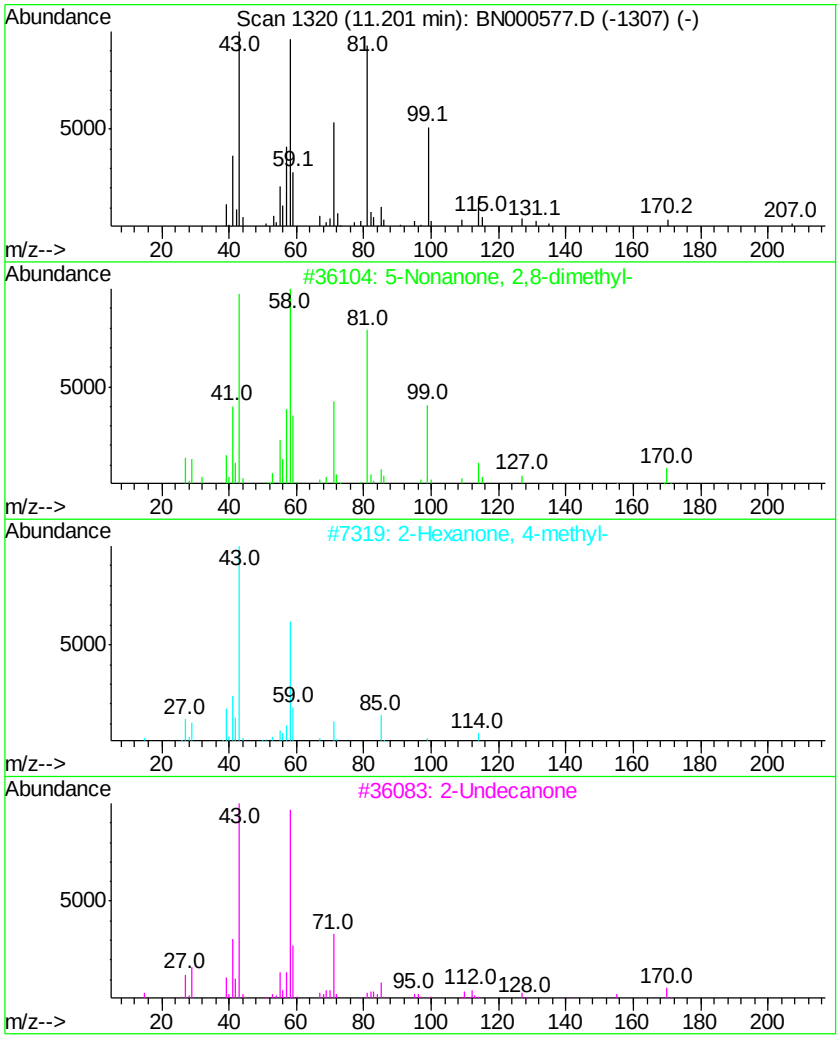
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

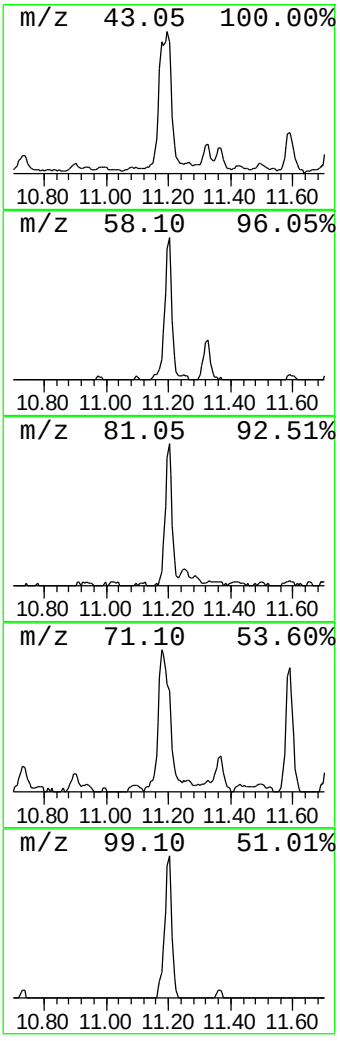
****

Peak Number 7 5-Nonanone, 2,8-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	4.38 ng/ul	343064	Naphthalene-d8	11.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Nonanone, 2,8-dimethyl-	170	C11H22O	002050-99-9	50
2	2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	35
3	2-Undecanone	170	C11H22O	000112-12-9	35
4	6-Undecanone	170	C11H22O	000927-49-1	32
5	6-Tridecanone	198	C13H26O	022026-12-6	32





Library Search Compound Report

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Operator : JU/SJ
Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

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TIC Library : C:\DATABASE\NIST02.L
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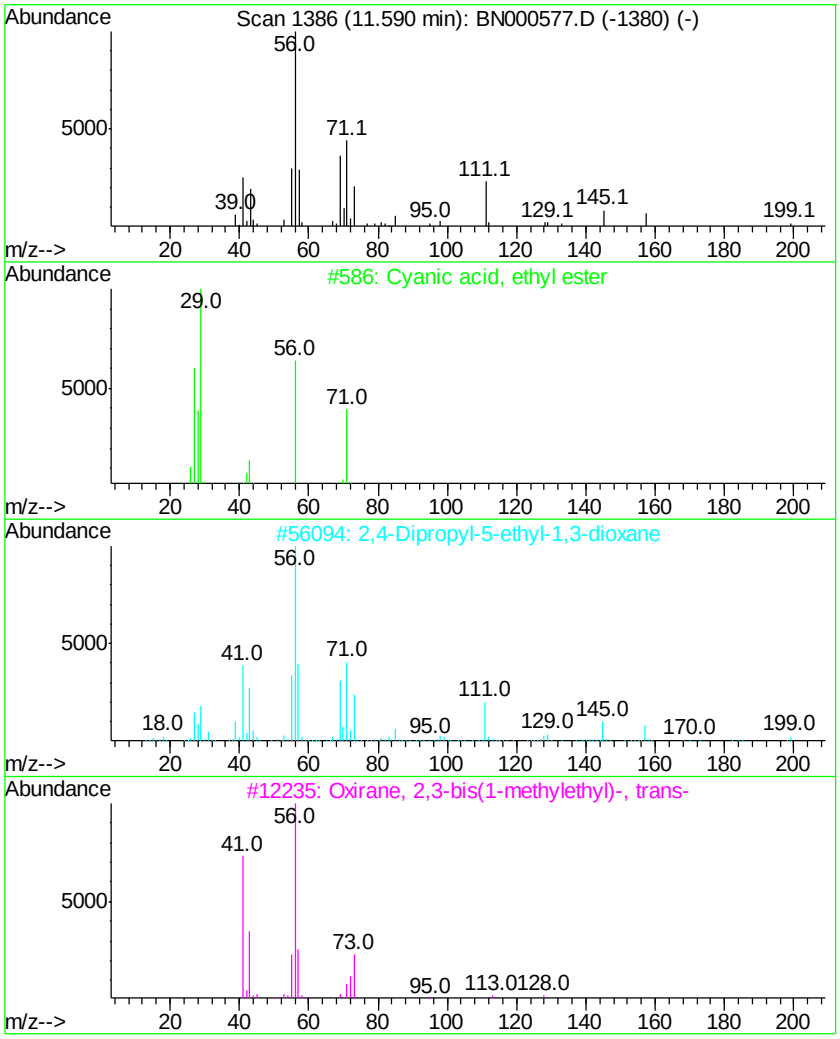
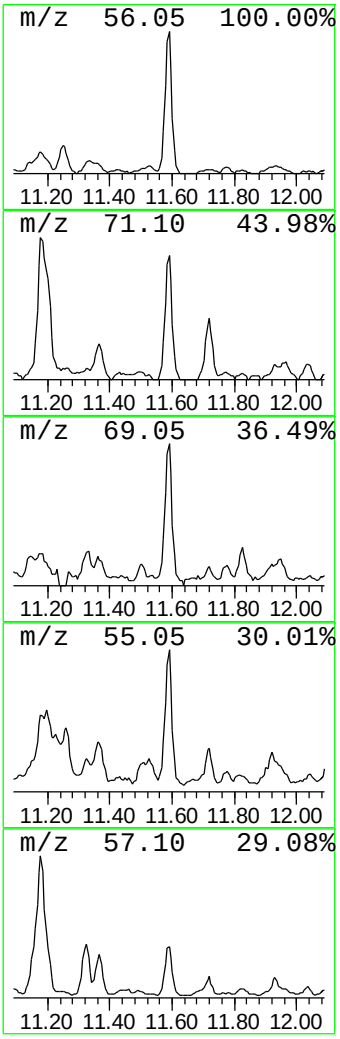
****

Peak Number 8 Cyanic acid, ethyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.59	2.15 ng/ul	168536	Naphthalene-d8	11.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	53
2	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	43
3	Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	37
4	2-Ethyl-cyclohexylamine	127	C8H17N	024216-90-8	35
5	1-Heptene, 2-methyl-	112	C8H16	015870-10-7	35





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Sample : J1905-07DL3 500X
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ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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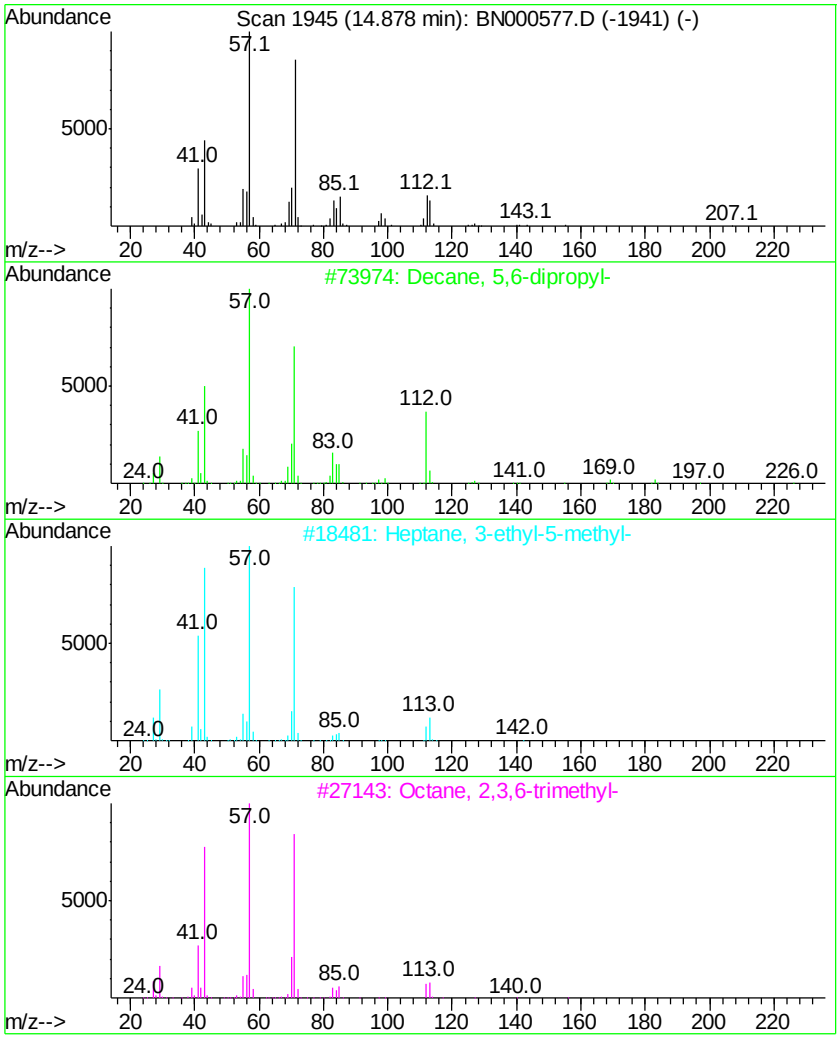
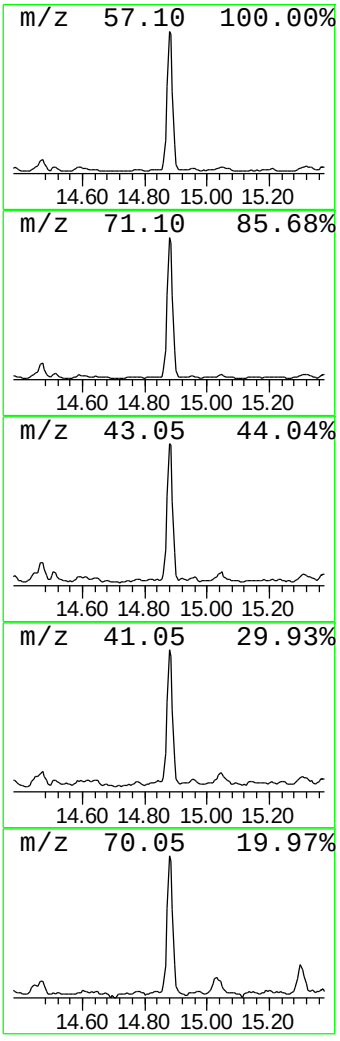
****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	4.26 ng/ul	446627	Acenaphthene-d10	15.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	72
2	Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	72
3	Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	72
4	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	59
5	Tridecane, 7-methyl-	198	C14H30	026730-14-3	59





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Sample : J1905-07DL3 500X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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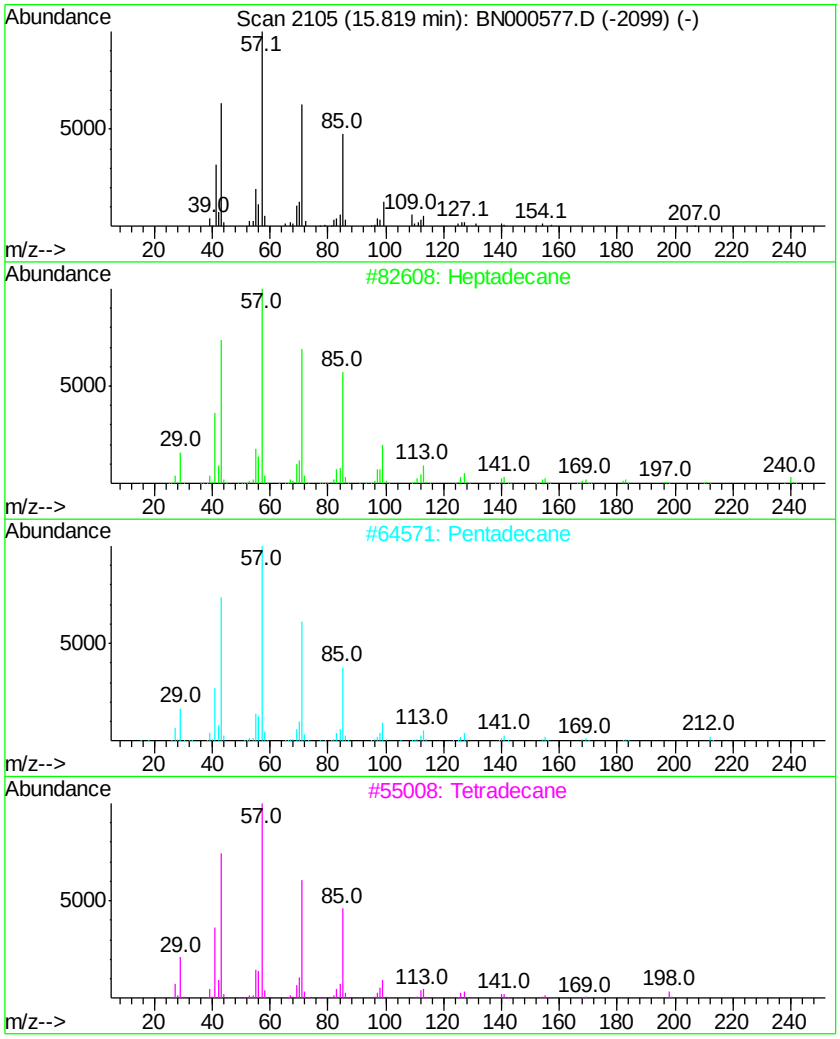
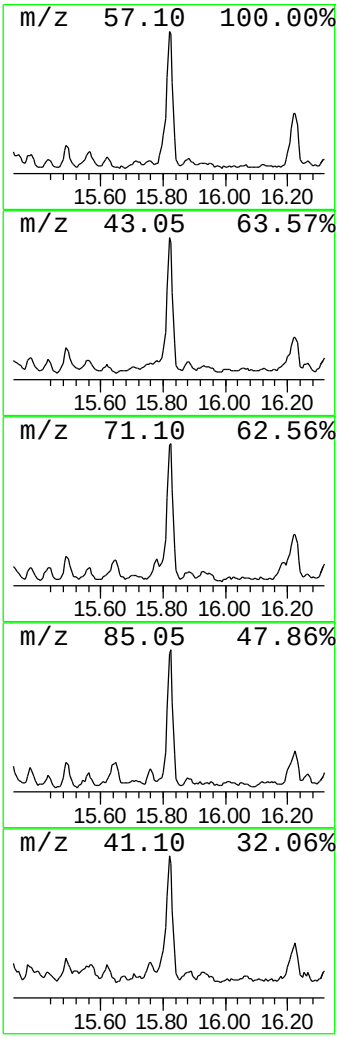
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	2.11 ng/ul	221304	Acenaphthene-d10	15.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	90
2	Pentadecane	212	C15H32	000629-62-9	86
3	Tetradecane	198	C14H30	000629-59-4	86
4	Hexadecane	226	C16H34	000544-76-3	83
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78





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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DAM42DL3

Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN031918.M
Quant Title : SVOA CALIBRATION

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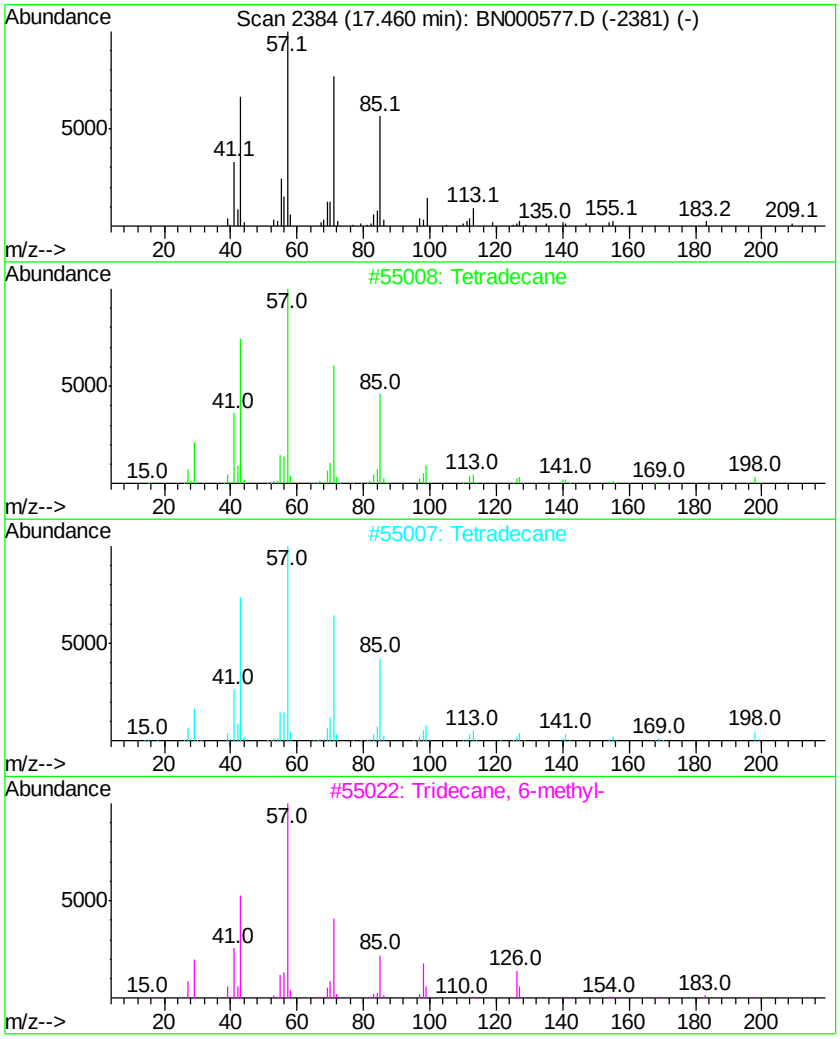
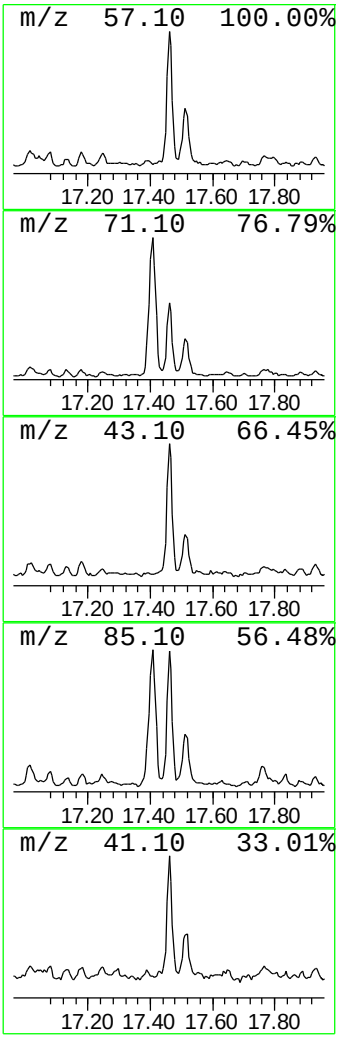
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.46	2.05 ng/ul	251608	Phenanthrene-d10	17.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	86
2	Tetradecane	198	C14H30	000629-59-4	86
3	Tridecane, 6-methyl-	198	C14H30	013287-21-3	72
4	Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	64





Tentatively Identified Compound (LSC) summary

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Heptanone, 4,6-...	7.80	2.0	ng/ul	106003	1	8.41	1055720	20.0
(DEL) Alkane: Str...	8.00	4.0	ng/ul	210856	1	8.41	1055720	20.0
1,1-Hexylenedioxy...	8.60	3.4	ng/ul	180517	1	8.41	1055720	20.0
Cyclohexanone, 3,...	8.84	3.5	ng/ul	183209	1	8.41	1055720	20.0
Benzene, 1,4-diet...	9.05	2.0	ng/ul	106908	1	8.41	1055720	20.0
(DEL) Alkane: Str...	9.63	5.4	ng/ul	283484	1	8.41	1055720	20.0
5-Nonanone, 2,8-d...	11.20	4.4	ng/ul	343064	2	11.25	1567190	20.0
Cyanic acid, ethy...	11.59	2.1	ng/ul	168536	2	11.25	1567190	20.0
(DEL) Alkane: Str...	14.88	4.3	ng/ul	446627	3	15.03	2097970	20.0
(DEL) Alkane: Str...	15.82	2.1	ng/ul	221304	3	15.03	2097970	20.0
(DEL) Alkane: Str...	17.46	2.0	ng/ul	251608	4	17.75	2456320	20.0