LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030220\
Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP002040.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.917	3	5	19	rVB	413042	580636	6.88%	0.590%
2	3.164	43	47	55	rBV	68461	94153	1.12%	0.096%
3	4.787	320	323	331	rVB	15344	25454	0.30%	0.026%
4	6.816	661	668	682	rBV	1167243	1899375	22.51%	1.929%
5	6.981	689	696	708	rVB	913815	1454302	17.23%	1.477%
6	7.175	721	729	742	rBV	1264646	2007918	23.79%	2.039%
7	7.640	799	808	816	rBV	1046330	1660205	19.67%	1.686%
8	8.346	920	928	940	rBV	1517452	2424199	28.72%	2.462%
9	8.781	994	1002	1016	rBV	1058065	1723420	20.42%	1.750%
10	9.269	1080	1085	1092	rVB	32368	48221	0.57%	0.049%
11	9.499	1116	1124	1133	rBV	881553	1422346	16.85%	1.445%
12	10.034	1207	1215	1240	rBV	1541625	2613944	30.97%	2.655%
13	10.410	1271	1279	1284	rBV	1477104	2434106	28.84%	2.472%
14	10.457	1284	1287	1294	rVB	345330	544352	6.45%	0.553%
15	10.546	1294	1302	1314	rBV	1440184	2521617	29.88%	2.561%
16	12.004	1541	1550	1556	rVV2	25722	55820	0.66%	0.057%
17	12.081	1556	1563	1573	rVV	281129	455437	5.40%	0.463%
18	12.204	1577	1584	1591	rVV7	12804	30356	0.36%	0.031%
19	12.304	1591	1601	1610	rVB	140936	243983	2.89%	0.248%
20	13.104	1731	1737	1746	rBV	133216	214577	2.54%	0.218%
21	13.193	1746	1752	1757	rVV2	17954	30495	0.36%	0.031%
22	13.304	1766	1771	1781	rVB2	26979	50516	0.60%	0.051%
23	13.445	1785	1795	1803	rBV4	29337	81436	0.96%	0.083%
24	13.598	1812	1821	1826	rBV2	47705	97213	1.15%	0.099%
25	13.651	1826	1830	1831	rVV2	36871	48526	0.57%	0.049%
26	13.687	1831	1836	1841	rVV	2634145	3696133	43.79%	3.754%
27	13.734	1841	1844	1850	rVV	309006	413307	4.90%	0.420%
28	13.834	1857	1861	1865	rBV2	19865	26732	0.32%	0.027%
29	13.963	1876	1883	1896	rBV	3231335	4797158	56.84%	4.872%
30	14.275	1928	1936	1942	rBV	2282055	3384083	40.10%	3.437%
31	14.340	1942	1947	1954	rVV	603113	881723	10.45%	0.896%
32	14.404	1954	1958	1964	rVB	67514	101671	1.20%	0.103%
33	14.475	1964	1970	1980	rBV	824856	1300377	15.41%	1.321%
34	14.551	1980	1983	1988	rVB	44661	65249	0.77%	0.066%

LSC Area Percent Report

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Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

35	14.675	1997	2004	2014	rVV	703539	1070486	12.68%	1.087%
36	14.763	2016	2019	2029	rVB7	11213	26247	0.31%	0.027%
37	15.075	2065	2072	2075	rBV2	20390	38031	0.45%	0.039%
38	15.169	2082	2088	2092	rVB5	15879	30721	0.36%	0.031%
39	15.275	2099	2106	2111	rBV	3997234	5831641	69.10%	5.923%
40	15.328	2111	2115	2120	rVB	458381	622384	7.37%	0.632%
41	15.392	2120	2126	2134	rBV	716572	997739	11.82%	1.013%
42	15.516	2140	2147	2150	rBV2	41373	91846	1.09%	0.093%
43	15.581	2155	2158	2161	rVB	23545	27940	0.33%	0.028%
44	15.628	2161	2166	2169	rBV	56348	83449	0.99%	0.085%
45	15.657	2169	2171	2178	rVB	58233	79623	0.94%	0.081%
46	15.775	2186	2191	2197	rVV	60804	114293	1.35%	0.116%
47	15.834	2197	2201	2204	rVV2	17066	27168	0.32%	0.028%
48	15.869	2204	2207	2214	rVB4	15323	26525	0.31%	0.027%
49	15.992	2222	2228	2232	rBV2	73554	133353	1.58%	0.135%
50	16.028	2232	2234	2238	rVV	48543	66490	0.79%	0.068%
51	16.063	2238	2240	2244	rVV3	22965	27224	0.32%	0.028%
52	16.104	2244	2247	2253	rVB	21709	28923	0.34%	0.029%
53	16.298	2275	2280	2284	rBV4	14324	28963	0.34%	0.029%
54	16.334	2284	2286	2291	rVV3	14954	24853	0.29%	0.025%
55	16.451	2301	2306	2315	rBV	47647	77442	0.92%	0.079%
56	16.681	2335	2345	2356	rVV2	298827	579737	6.87%	0.589%
57	16.763	2356	2359	2364	rVV4	9924	21910	0.26%	0.022%
58	16.845	2364	2373	2378	rVB	167222	263910	3.13%	0.268%
59	16.986	2392	2397	2398	rBV	25519	35287	0.42%	0.036%
60	17.028	2398	2404	2407	rVV	2868776	3978748	47.14%	4.041%
61	17.069	2407	2411	2415	rVV	2355758	3319793	39.34%	3.372%
62	17.128	2415	2421	2433	rVB	5452345	8032399	95.17%	8.158%
63	17.351	2454	2459	2462	rBV3	13834	24535	0.29%	0.025%
64	17.428	2462	2472	2479	rVV	243005	399924	4.74%	0.406%
65	17.569	2490	2496	2501	rBV	164401	246782	2.92%	0.251%
66	17.904	2548	2553	2556	rBV	73952	107160	1.27%	0.109%
67	17.945	2556	2560	2567	rBV	304303	421853	5.00%	0.428%
68	18.092	2579	2585	2588	rBV2	103033	163073	1.93%	0.166%
69	18.128	2588	2591	2598	rVB4	50255	73491	0.87%	0.075%
70	18.198	2598	2603	2606	rBV2	21852	43805	0.52%	0.044%
71	18.233	2606	2609	2618	rVV4	59654	94911	1.12%	0.096%

LSC Area Percent Report

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Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

72	18.328	2622	2625	2628	rVV2	18544	25770	0.31%	0.026%
73	18.392	2629	2636	2638	rVV	68224	121187	1.44%	0.123%
74	18.422	2638	2641	2649	rVB2	106382	159368	1.89%	0.162%
75	18.610	2670	2673	2680	rVB4	46054	72660	0.86%	0.074%
76	18.757	2694	2698	2701	rVV	24591	35767	0.42%	0.036%
77	18.792	2701	2704	2709	rVB4	16405	27937	0.33%	0.028%
78	18.928	2722	2727	2734	rBV	35680	60192	0.71%	0.061%
79	19.045	2738	2747	2754	rBV	924183	1249891	14.81%	1.269%
80	19.186	2768	2771	2776	rBV5	32531	45701	0.54%	0.046%
81	19.263	2779	2784	2793	rVB2	48992	96231	1.14%	0.098%
82	19.369	2796	2802	2814	rBV2	5438705	8439650	100.00%	8.572%
83	19.575	2834	2837	2842	rVB	25755	27612	0.33%	0.028%
84	19.886	2887	2890	2894	rVB	29226	37920	0.45%	0.039%
85	19.980	2903	2906	2911	rVB	30613	37920	0.45%	0.039%
86	20.469	2986	2989	2995	rVB2	28861	39996	0.47%	0.041%
87	20.863	3051	3056	3073	rBV4	413582	796856	9.44%	0.809%
88	21.122	3093	3100	3105	rBV	3713634	4916667	58.26%	4.994%
89	21.233	3116	3119	3123	rBV2	41956	66323	0.79%	0.067%
90	21.298	3127	3130	3137	rVB	166679	181716	2.15%	0.185%
91	21.727	3199	3203	3216	rVB2	315836	434608	5.15%	0.441%
92	22.186	3277	3281	3287	rBV	416275	525137	6.22%	0.533%
93	22.316	3298	3303	3310	rBV	294155	389091	4.61%	0.395%
94	22.692	3362	3367	3377	rVB	488187	761980	9.03%	0.774%
95	23.186	3444	3451	3459	rBV	4277976	7956231	94.27%	8.081%
96	23.263	3459	3464	3468	rVV	485075	799946	9.48%	0.812%
97	23.327	3468	3475	3492	rVB	2645755	5022136	59.51%	5.101%
98	23.910	3568	3574	3582	rBV	364408	692036	8.20%	0.703%
99	24.657	3695	3701	3711	rVB	227710	485762	5.76%	0.493%
100	25.533	3845	3850	3861	rVB3	102395	262009	3.10%	0.266%

Sum of corrected areas: 98459999

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030220\
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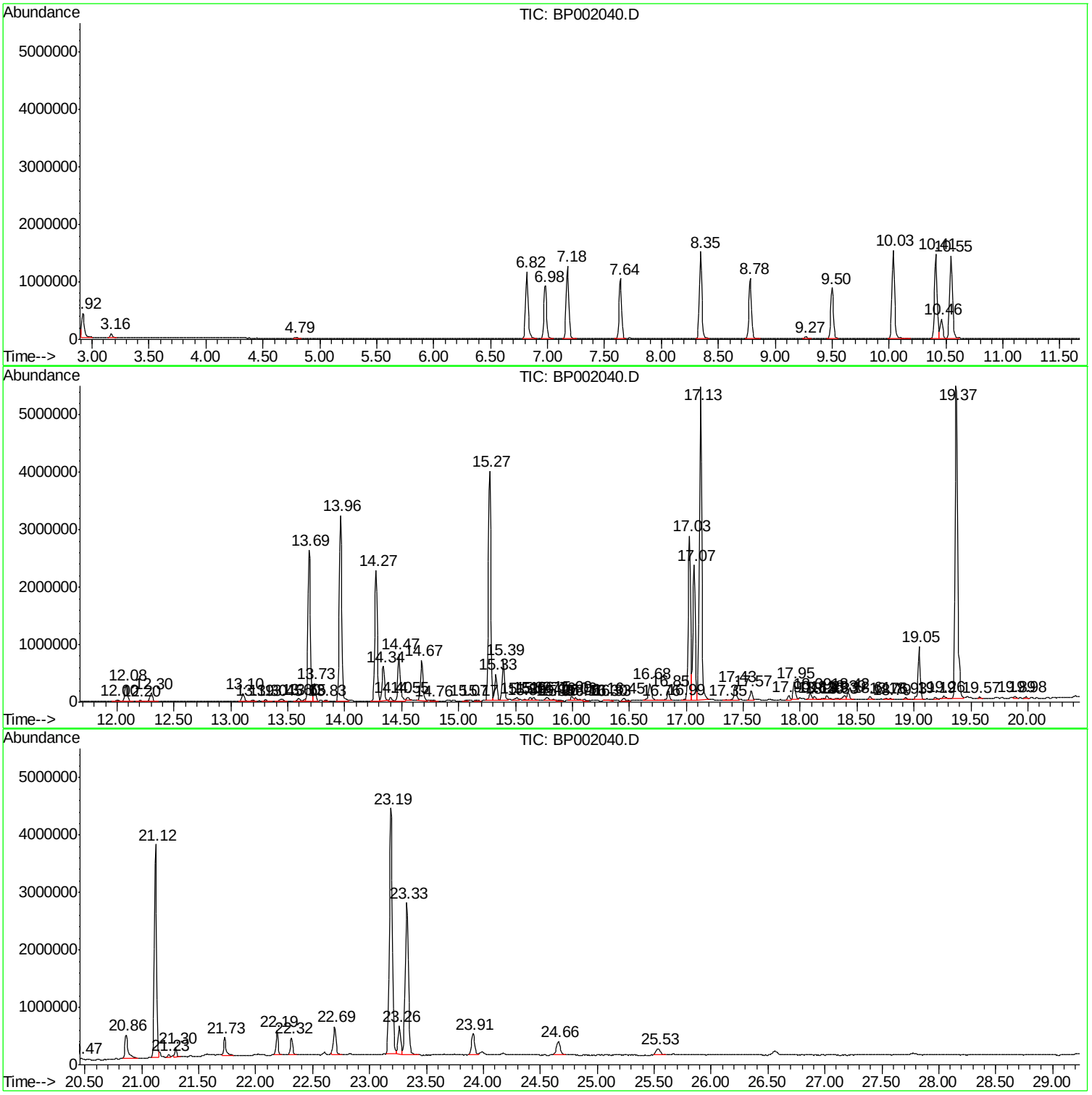
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

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Quant Title : SVOA CALIBRATION

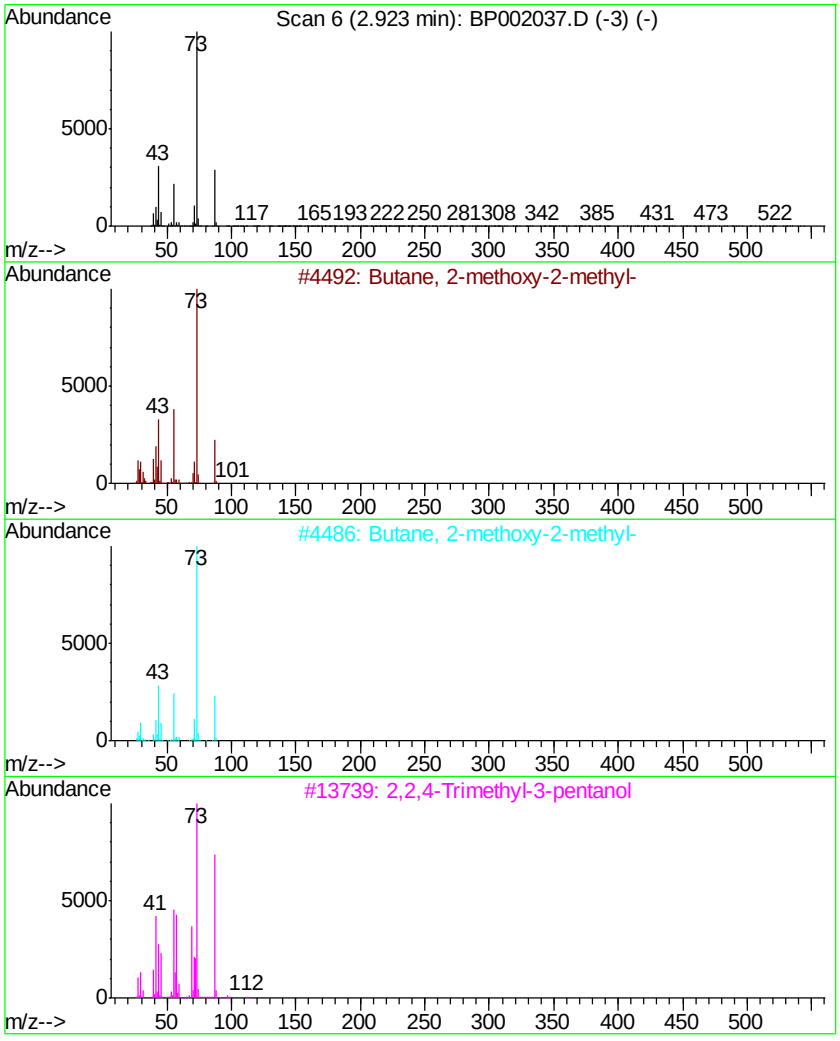
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

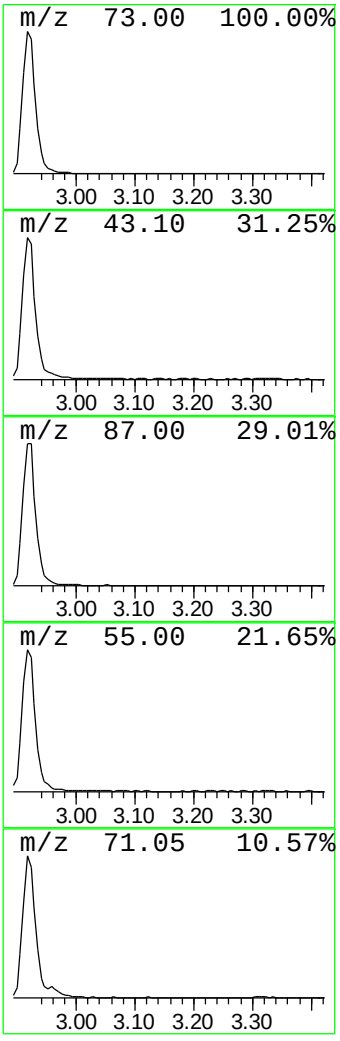
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	6.99 ng/ul	580636	1,4-Dichlorobenzene-d4	7.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030220\
Data File : BP002037.D
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Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

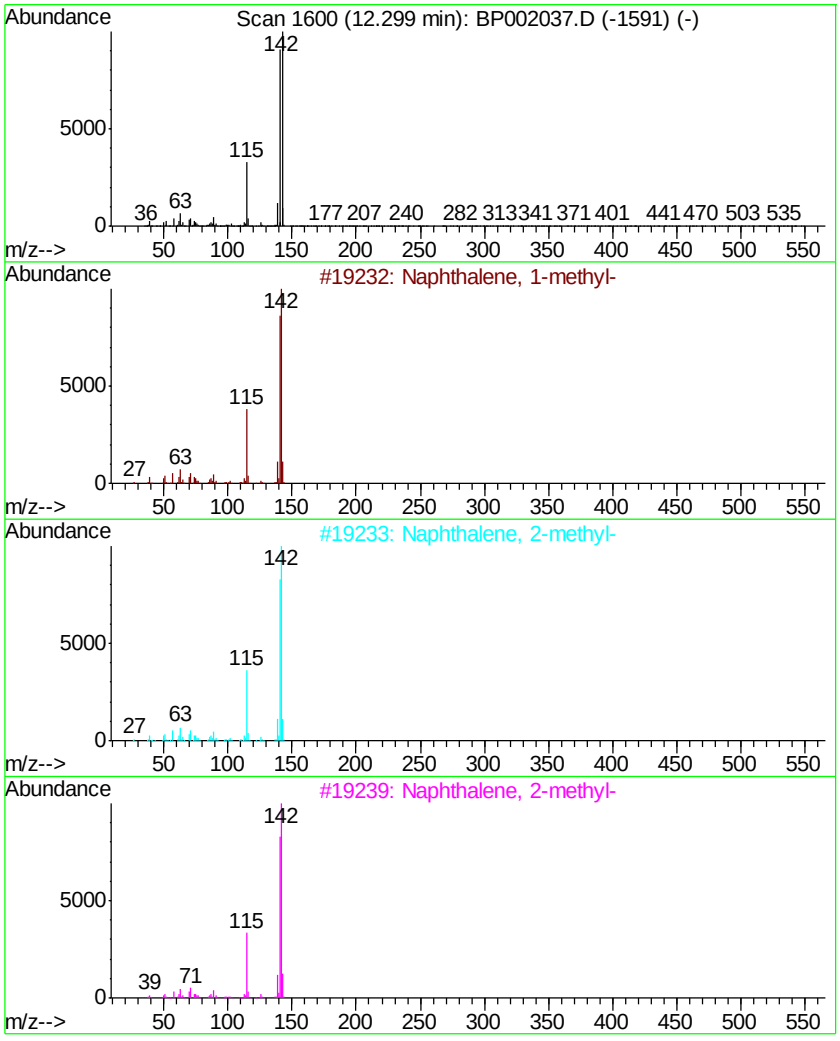
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

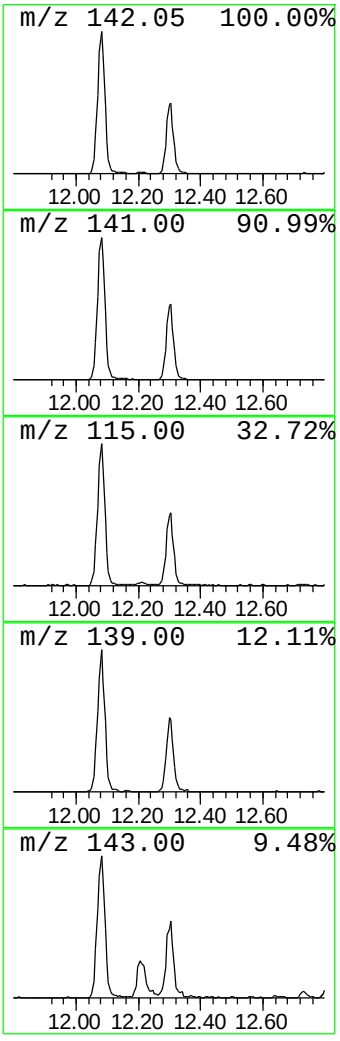
****

Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	2.00 ng/ul	243983	Naphthalene-d8	10.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5	Benzocycloheptatriene	142	C11H10	000264-09-5	91





Library Search Compound Report

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ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

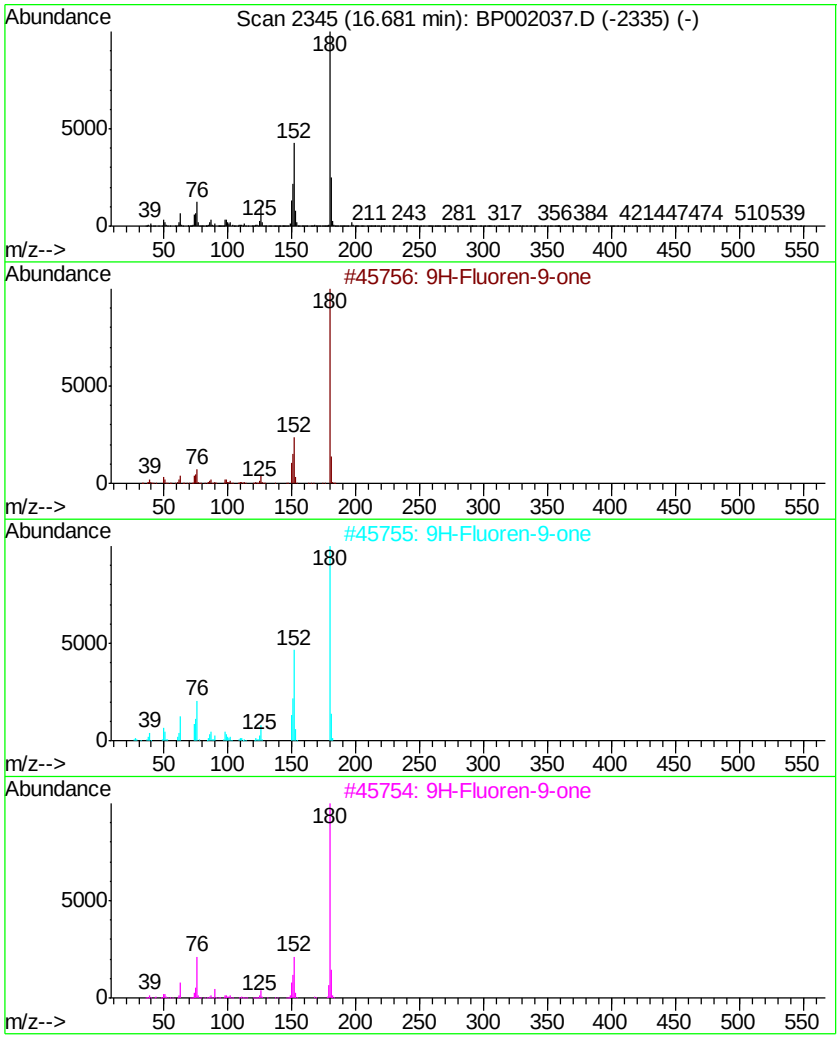
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

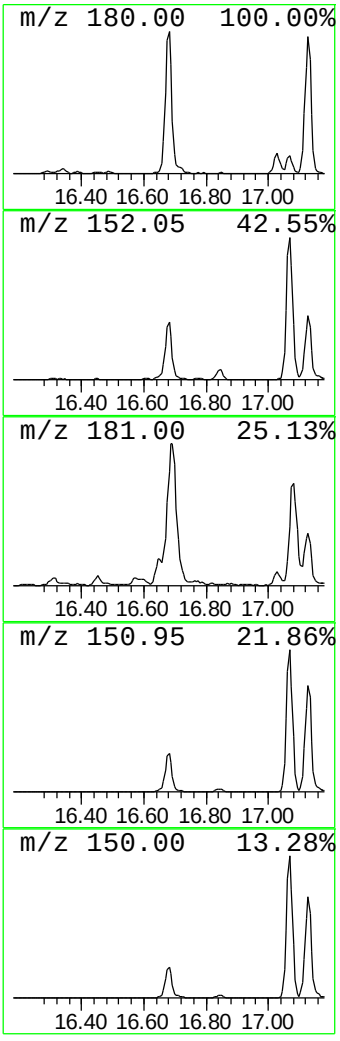
****

Peak Number 3 9H-Fluoren-9-one Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.68	2.91 ng/ul	579737	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one	180	C13H8O	000486-25-9	93
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	93
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	90
4	9H-Fluoren-9-one	180	C13H8O	000486-25-9	90
5	Benzo[h]cinnoline	180	C12H8N2	000230-31-9	64





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Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

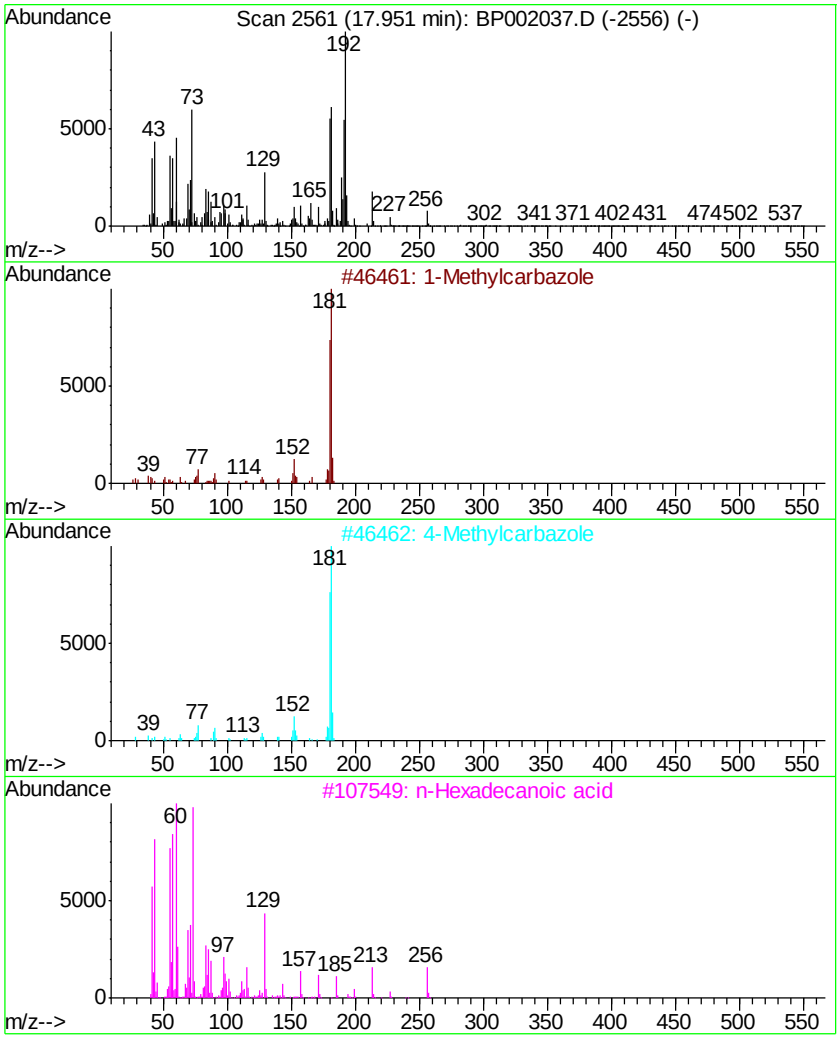
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

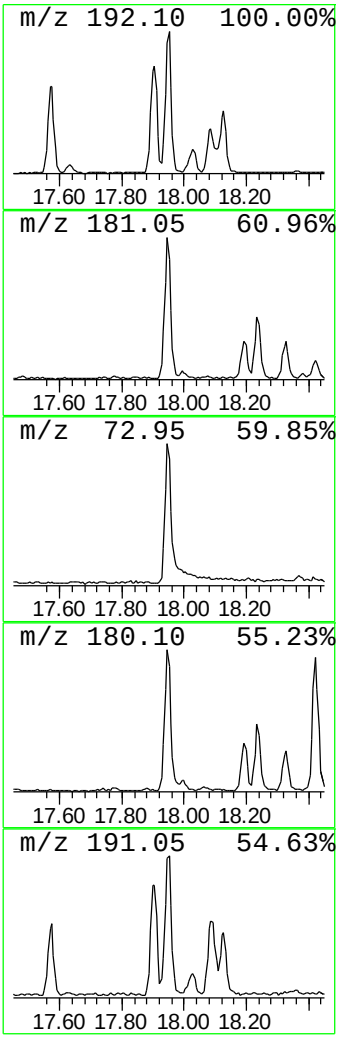
****

Peak Number 4 1-Methylcarbazole Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.95	2.12 ng/ul	421853	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methylcarbazole	181	C13H11N	006510-65-2	91
2	4-Methylcarbazole	181	C13H11N	003770-48-7	91
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
4	3-Methylcarbazole	181	C13H11N	004630-20-0	68
5	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	68





Library Search Compound Report

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Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

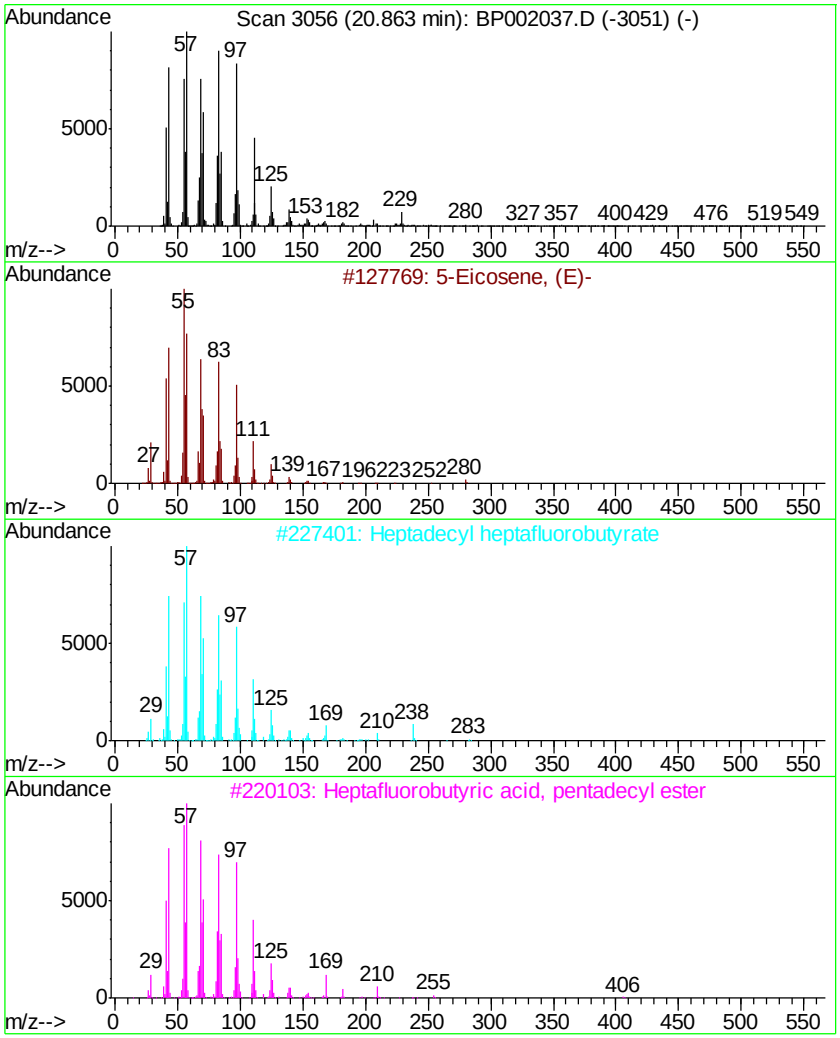
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

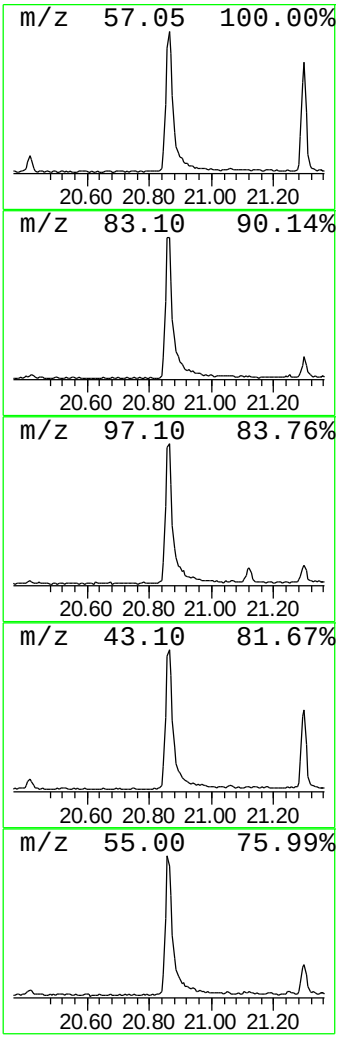
****

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	3.24 ng/ul	796856	Chrysene-d12	21.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4	1-Triacontanol	438	C30H62O	000593-50-0	93
5	Cyclopentadecane	210	C15H30	000295-48-7	93





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Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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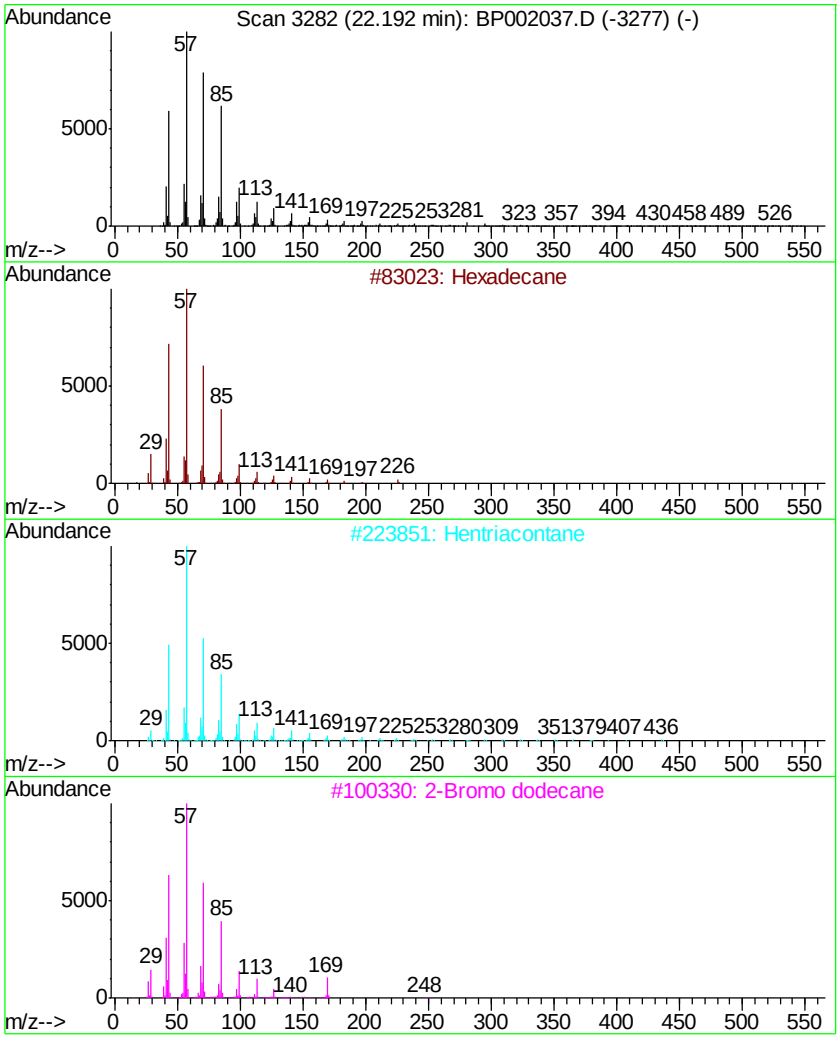
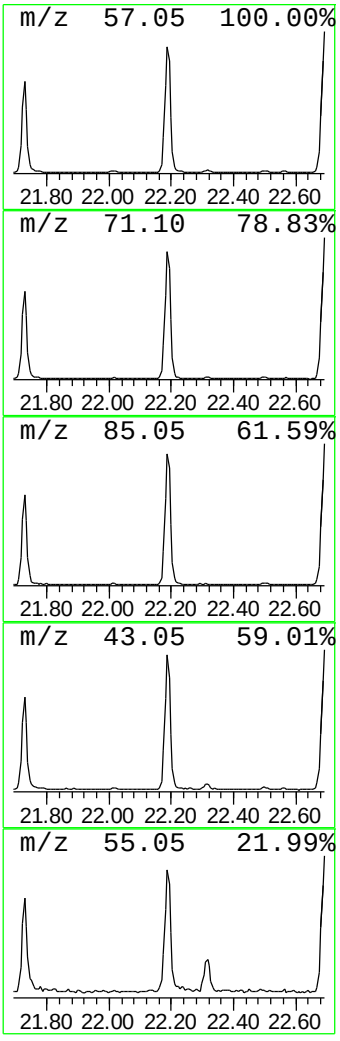
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.19	2.14 ng/ul	525137	Chrysene-d12	21.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	95
2	Hentriacontane	437	C31H64	000630-04-6	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
5	Heneicosane	296	C21H44	000629-94-7	91





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Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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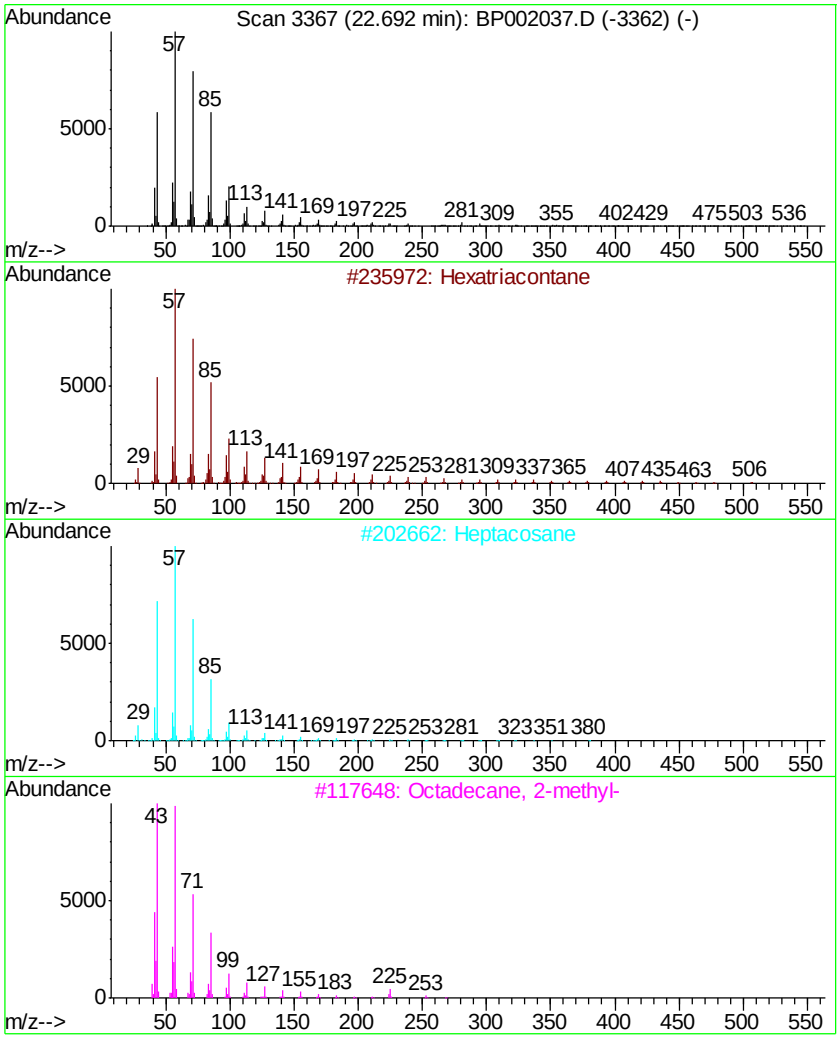
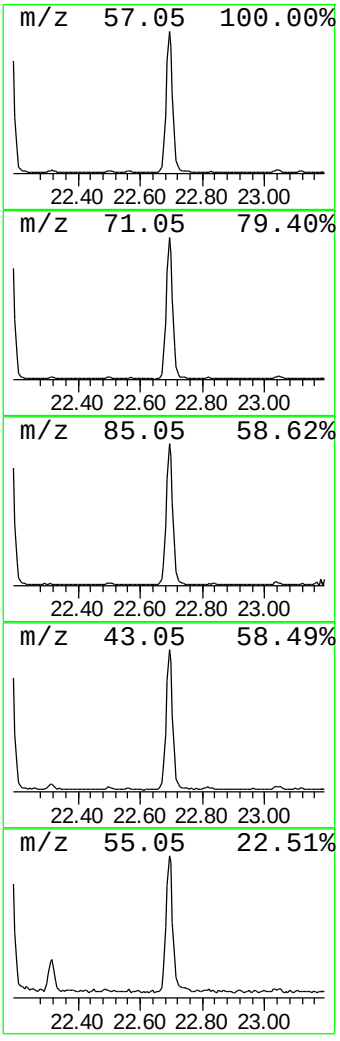
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.69	3.03 ng/ul	761980	Perylene-d12	23.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	93
2	Heptacosane	380	C27H56	000593-49-7	91
3	Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Triacontane	422	C30H62	000638-68-6	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030220\
Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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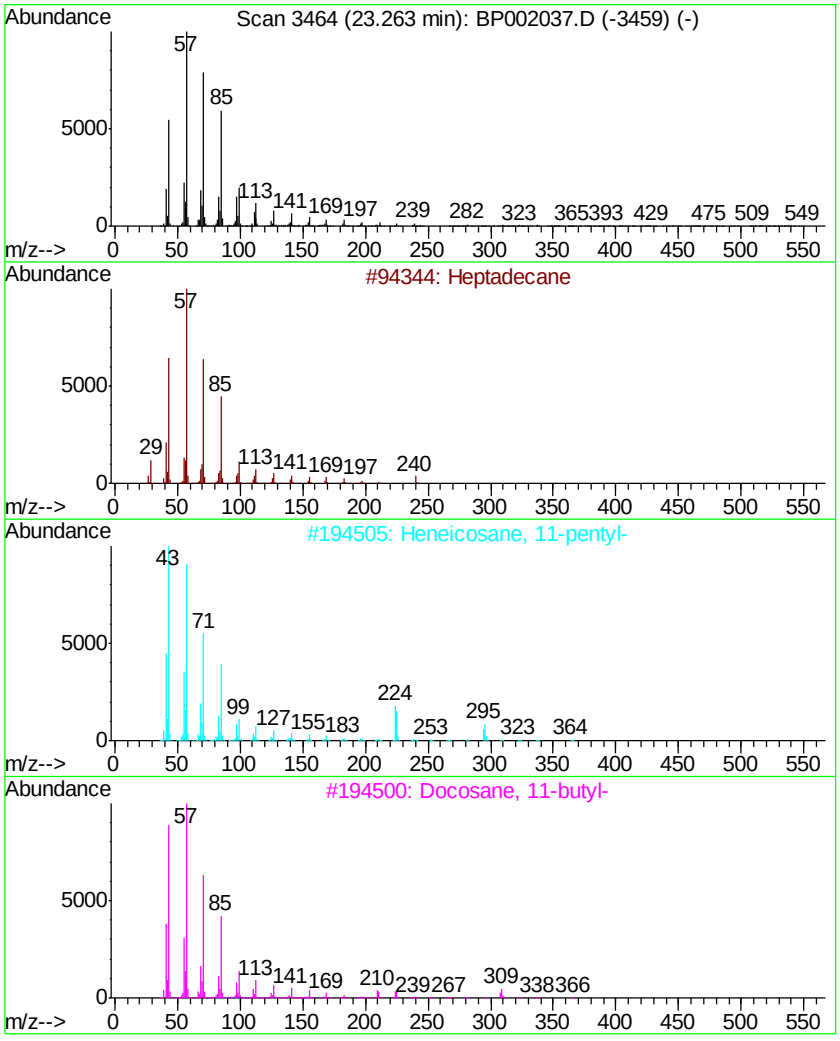
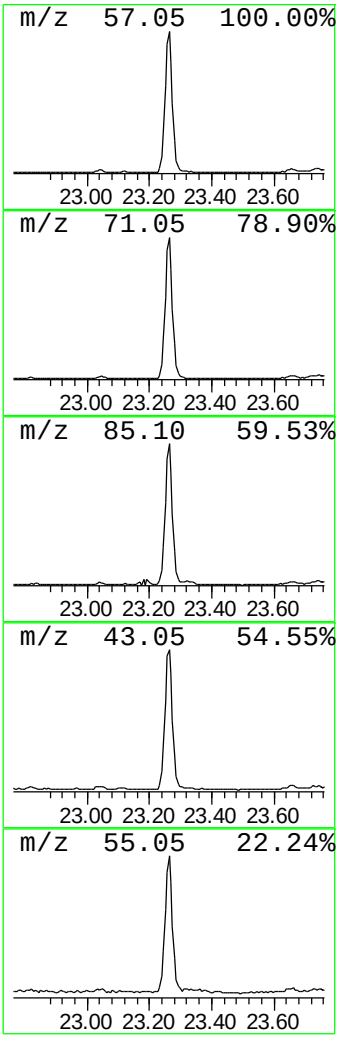
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.26	3.19 ng/ul	799946	Perylene-d12	23.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	94
2	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
3	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
4	Tetracosane	338	C24H50	000646-31-1	91
5	Tetratetracontane	619	C44H90	007098-22-8	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030220\
Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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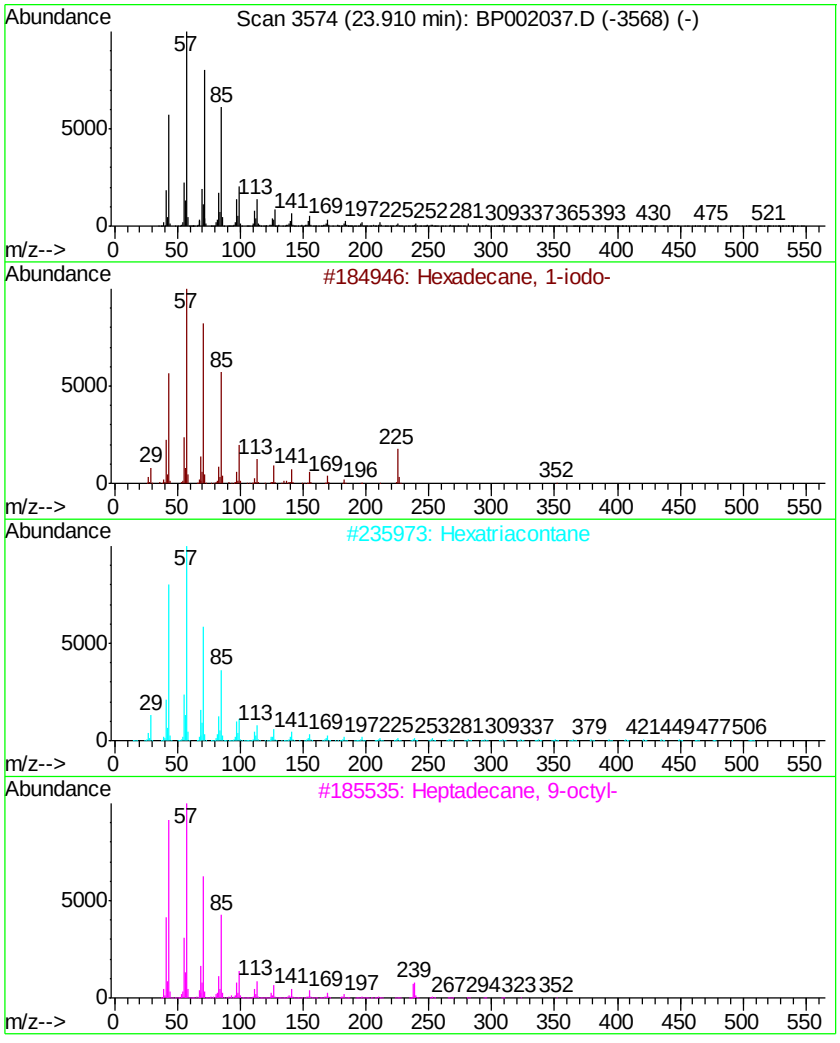
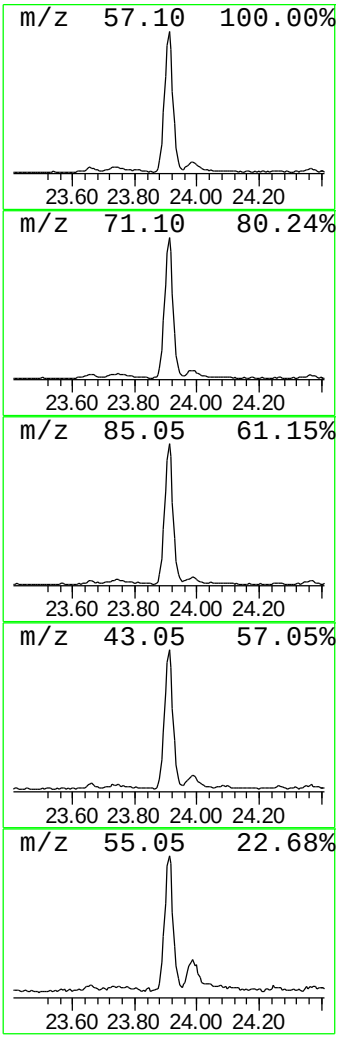
Peak Number 9 Hexadecane, 1-iodo- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.91	2.76 ng/ul	692036	Perylene-d12	23.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
2	Hexatriacontane	507	C36H74	000630-06-8	95
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4	Hexatriacontane	507	C36H74	000630-06-8	93
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030220\
Data File : BP002037.D
Acq On : 04 Mar 2020 02:22
Operator : CG/JU
Sample : L1617-21
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBDT5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	2.92	7.0	ng/ul	580636	1	7.64	1660210	20.0
Naphthalene, 1-me...	12.30	2.0	ng/ul	243983	2	10.41	2434110	20.0
9H-Fluoren-9-one	16.68	2.9	ng/ul	579737	4	17.03	3978750	20.0
1-Methylcarbazole	17.95	2.1	ng/ul	421853	4	17.03	3978750	20.0
(DEL) Alkane: Str...	20.86	3.2	ng/ul	796856	5	21.12	4916670	20.0
(DEL) Alkane: Str...	22.19	2.1	ng/ul	525137	5	21.12	4916670	20.0
(DEL) Alkane: Str...	22.69	3.0	ng/ul	761980	6	23.33	5022140	20.0
(DEL) Alkane: Str...	23.26	3.2	ng/ul	799946	6	23.33	5022140	20.0
Hexadecane, 1-iodo-	23.91	2.8	ng/ul	692036	6	23.33	5022140	20.0