LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

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_M\METHODS\SOM-EPA-BM070819MA.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	4.369	232	235	241	rVB	90307	108635	3.73%	0.277%
2	4.857	315	318	323	rVB	47395	61324	2.10%	0.156%
3	6.393	576	579	585	rVB2	55517	80392	2.76%	0.205%
4	6.816	648	651	655	rVB	44848	57152	1.96%	0.146%
5	6.893	659	664	678	rVB	387378	658249	22.57%	1.680%
6	7.063	688	693	702	rVB	287403	457324	15.68%	1.167%
7	7.251	720	725	735	rVB	410502	665048	22.80%	1.697%
8	7.716	799	804	812	rVB	496928	791605	27.14%	2.020%
9	8.428	918	925	938	rBV	422592	712937	24.45%	1.819%
10	8.881	996	1002	1011	rBV	378566	637242	21.85%	1.626%
11	9.240	1058	1063	1071	rVB	22906	31448	1.08%	0.080%
12	9.598	1118	1124	1133	rBV	296520	482214	16.54%	1.231%
13	9.816	1155	1161	1167	rVB5	6392	13931	0.48%	0.036%
14	10.134	1208	1215	1229	rVB	446116	843875	28.94%	2.153%
15	10.504	1271	1278	1285	rBV	632825	1063134	36.46%	2.713%
16	10.669	1299	1306	1317	rBV	82224	169437	5.81%	0.432%
17	11.728	1482	1486	1490	rVB	13813	17405	0.60%	0.044%
18	12.104	1546	1550	1555	rVB3	5522	10006	0.34%	0.026%
19	12.263	1571	1577	1585	rBV2	14223	29139	1.00%	0.074%
20	12.698	1647	1651	1656	rVB3	14185	20658	0.71%	0.053%
21	12.957	1691	1695	1699	rBV3	6387	8411	0.29%	0.021%
22	13.222	1735	1740	1744	rVB3	16131	22492	0.77%	0.057%
23	13.669	1812	1816	1820	rVB3	14312	17569	0.60%	0.045%
24	13.780	1825	1835	1840	rVV	820226	1221258	41.88%	3.116%
25	13.827	1840	1843	1851	rVB	85742	120517	4.13%	0.308%
26	14.057	1871	1882	1888	rBV	998255	1458153	50.00%	3.721%
27	14.104	1889	1890	1896	rVB3	12713	13630	0.47%	0.035%
28	14.192	1900	1905	1909	rVV2	24043	29722	1.02%	0.076%
29	14.310	1920	1925	1928	rBV	85746	113751	3.90%	0.290%
30	14.363	1928	1934	1940	rVB2	959779	1417233	48.60%	3.617%
31	14.416	1940	1943	1946	rBV2	13653	17854	0.61%	0.046%
32	14.592	1967	1973	1984	rBV2	184625	393873	13.51%	1.005%
33	14.674	1984	1987	1992	rVB2	15530	22693	0.78%	0.058%
34	14.769	1998	2003	2004	rBV2	14227	21148	0.73%	0.054%

LSC Area Percent Report

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Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

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_M\METHODS\SOM-EPA-BM070819MA.M
Title : SVOA CALIBRATION

35	15.157	2063	2069	2073	rBV2	8073	10398	0.36%	0.027%
36	15.192	2073	2075	2081	rVB4	4980	8306	0.28%	0.021%
37	15.357	2097	2103	2114	rBV	1262394	1865059	63.95%	4.759%
38	15.498	2121	2127	2132	rBV	212493	324834	11.14%	0.829%
39	15.545	2132	2135	2140	rVV3	35627	48564	1.67%	0.124%
40	15.604	2140	2145	2149	rVB3	14428	19933	0.68%	0.051%
41	15.739	2164	2168	2171	rBV4	15538	22584	0.77%	0.058%
42	15.774	2171	2174	2178	rVV3	18746	25130	0.86%	0.064%
43	15.833	2179	2184	2191	rVB3	15649	30709	1.05%	0.078%
44	15.892	2191	2194	2197	rBV3	5913	8720	0.30%	0.022%
45	15.927	2197	2200	2204	rVB5	9415	11750	0.40%	0.030%
46	16.069	2221	2224	2231	rBV8	7437	15808	0.54%	0.040%
47	16.210	2244	2248	2250	rBV4	14106	17670	0.61%	0.045%
48	16.521	2296	2301	2305	rBV4	7202	10478	0.36%	0.027%
49	17.010	2380	2384	2391	rBV2	56083	87428	3.00%	0.223%
50	17.080	2391	2396	2397	rVV2	40615	56708	1.94%	0.145%
51	17.115	2397	2402	2411	rVV2	1235197	1719015	58.95%	4.387%
52	17.215	2413	2419	2428	rVV	1216311	1764479	60.51%	4.503%
53	17.398	2447	2450	2456	rVB4	9157	13057	0.45%	0.033%
54	17.598	2482	2484	2493	rBV7	6346	14050	0.48%	0.036%
55	17.745	2506	2509	2513	rVB6	8810	13019	0.45%	0.033%
56	17.892	2528	2534	2538	rBV6	15011	26996	0.93%	0.069%
57	17.951	2540	2544	2549	rVV2	27424	45759	1.57%	0.117%
58	18.010	2549	2554	2564	rVV	271311	455549	15.62%	1.162%
59	18.080	2564	2566	2574	rVB	19275	37723	1.29%	0.096%
60	18.310	2602	2605	2609	rBV6	11410	17881	0.61%	0.046%
61	18.351	2610	2612	2616	rVB5	12204	14748	0.51%	0.038%
62	18.521	2636	2641	2646	rBV8	11582	22238	0.76%	0.057%
63	18.686	2666	2669	2672	rBV5	7772	8389	0.29%	0.021%
64	18.927	2707	2710	2715	rVB5	25672	31279	1.07%	0.080%
65	19.145	2743	2747	2749	rBV2	45908	57036	1.96%	0.146%
66	19.174	2749	2752	2754	rVV3	17635	25658	0.88%	0.065%
67	19.292	2768	2772	2783	rBV	183081	301439	10.34%	0.769%
68	19.509	2803	2809	2821	rBV	1687450	2406773	82.53%	6.142%
69	19.904	2873	2876	2881	rVB	52933	63503	2.18%	0.162%
70	20.051	2898	2901	2904	rBV3	62535	62480	2.14%	0.159%
71	20.551	2982	2986	2991	rBV7	49029	81617	2.80%	0.208%

LSC Area Percent Report

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Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

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_M\METHODS\SOM-EPA-BM070819MA.M
Title : SVOA CALIBRATION

72	20.862	3034	3039	3044	rBV	1260934	1382340	47.40%	3.528%
73	20.986	3057	3060	3062	rBV	123926	161068	5.52%	0.411%
74	21.015	3062	3065	3075	rVB	332092	538848	18.48%	1.375%
75	21.303	3109	3114	3123	rBV	1672202	2181965	74.82%	5.568%
76	21.880	3209	3212	3214	rBV	131119	157874	5.41%	0.403%
77	21.909	3214	3217	3227	rVB2	380490	609148	20.89%	1.554%
78	22.850	3372	3377	3382	rBV	1088298	1527457	52.38%	3.898%
79	23.415	3467	3473	3482	rVB	1659392	2916239	100.00%	7.442%
80	23.556	3491	3497	3508	rVB2	1177925	2274602	78.00%	5.804%
81	24.062	3577	3583	3591	rVB	1336832	2562432	87.87%	6.539%
82	24.156	3594	3599	3612	rVB2	573456	1456706	49.95%	3.717%
83	26.603	4008	4015	4030	rVB5	560890	1882358	64.55%	4.803%

Sum of corrected areas: 39187263

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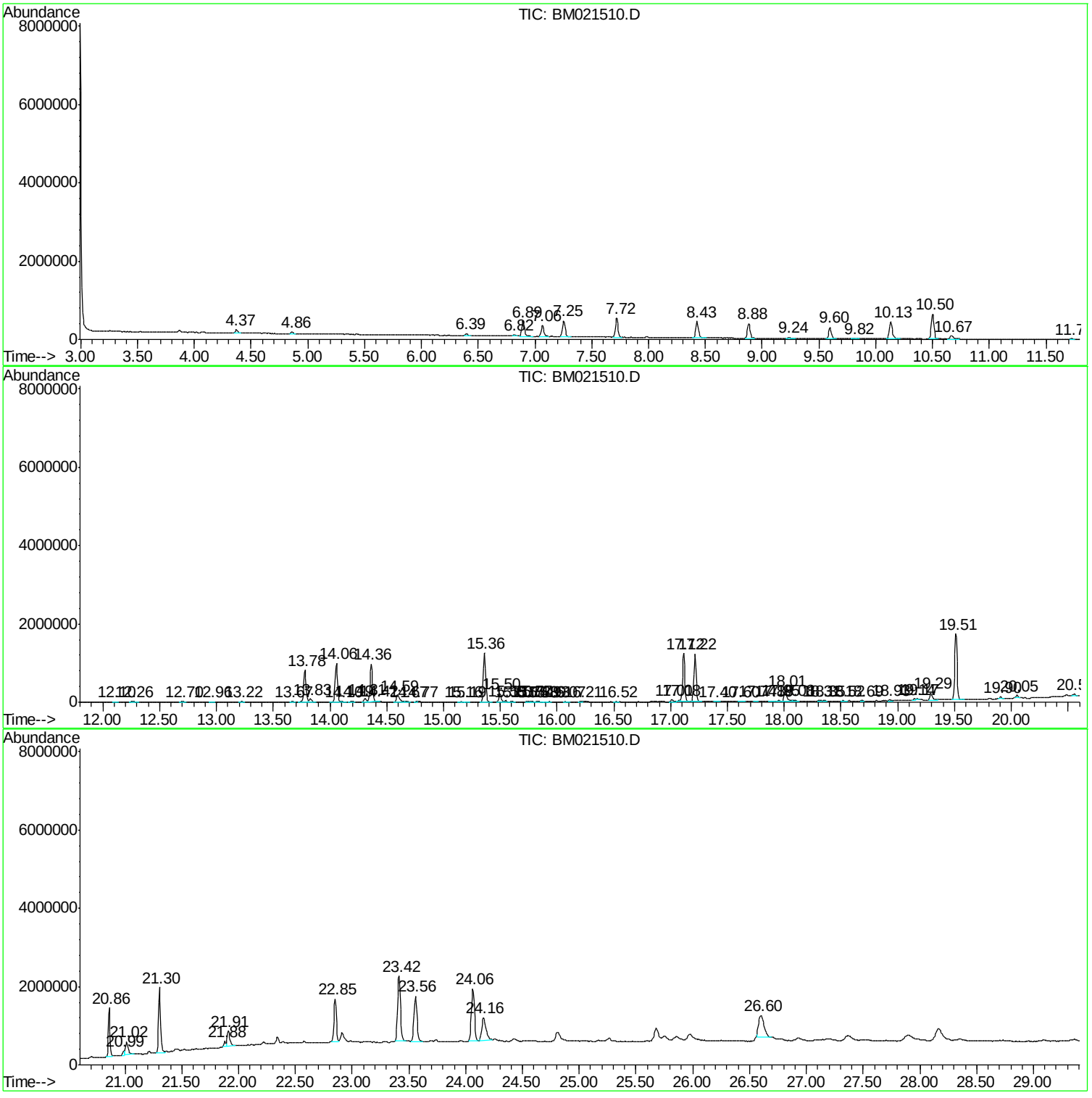
Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
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Misc :
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Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

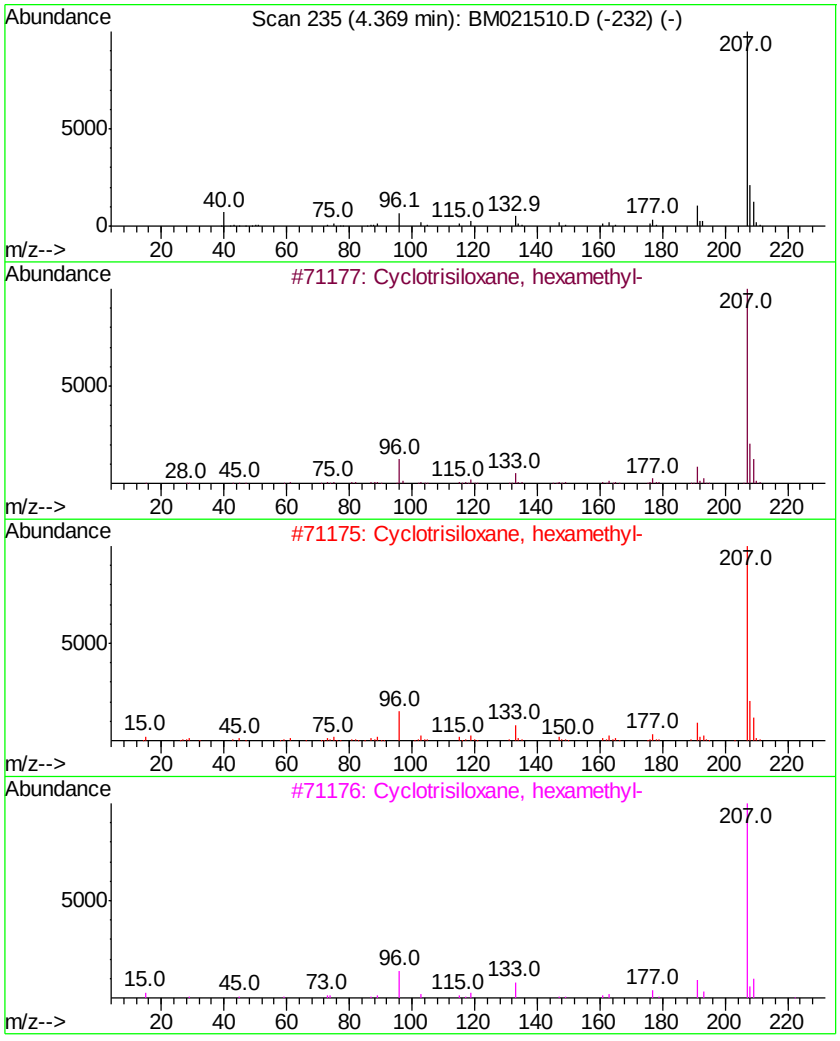
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

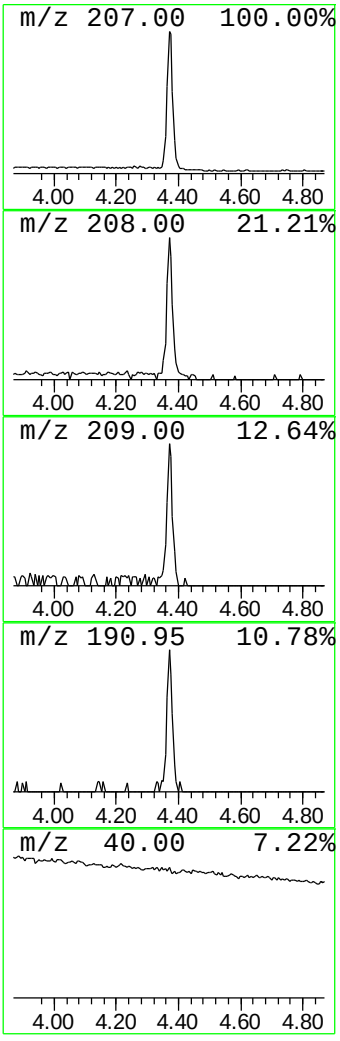
****

Peak Number 1 Cyclotrisiloxane, hexamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	2.74 ng/ul	108635	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	90
2	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	86
3	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	72
4	Benzo[h]quinoline, 2,4-dimethyl-	207	C15H13N	000605-67-4	39
5	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	39





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
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Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

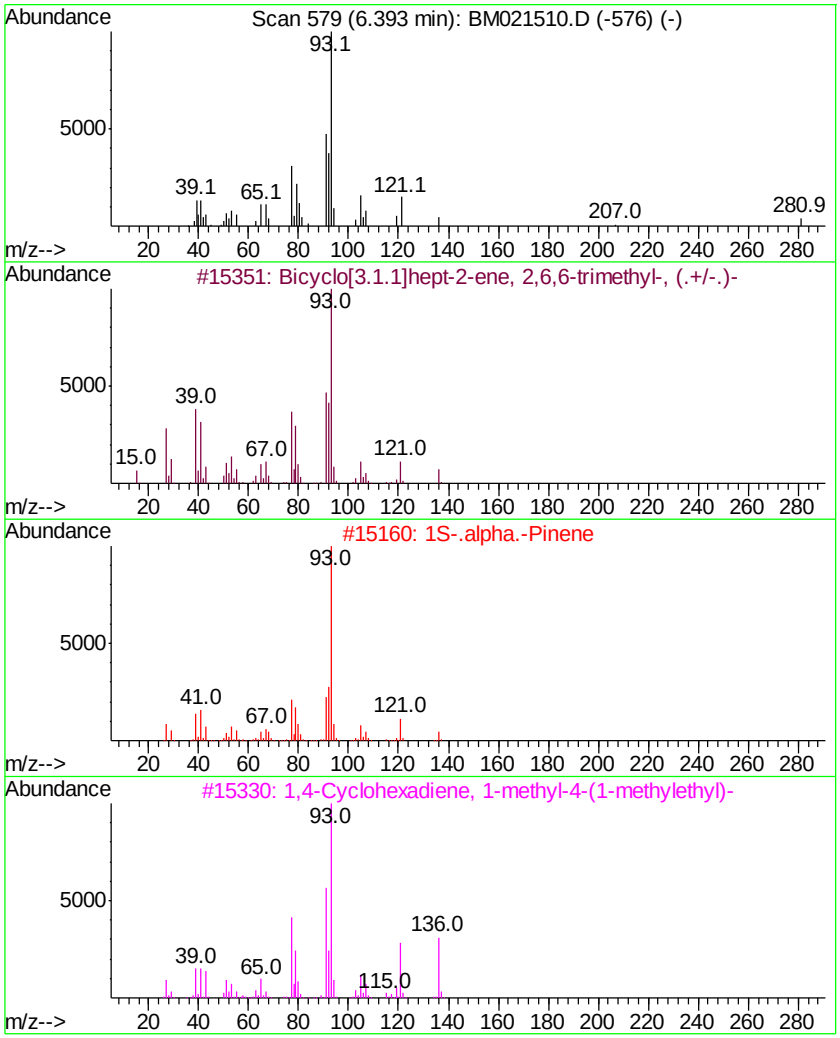
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

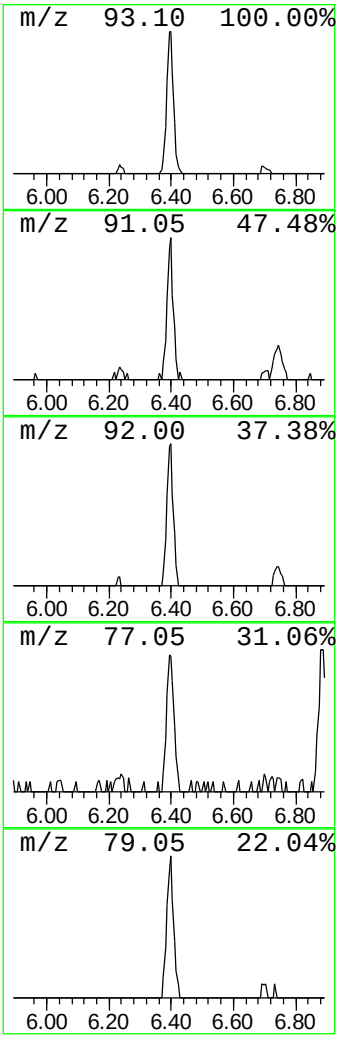
****

Peak Number 2 Bicyclo[3.1.1]hept-2-ene, 2... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	2.03 ng/ul	80392	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.1]hept-2-ene, 2,6,6-...	136	C10H16	002437-95-8	94
2	1S-.alpha.-Pinene	136	C10H16	007785-26-4	91
3	1,4-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-85-4	91
4	1,4-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-85-4	91
5	.alpha.-Pinene	136	C10H16	000080-56-8	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
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Instrument : BNA_M
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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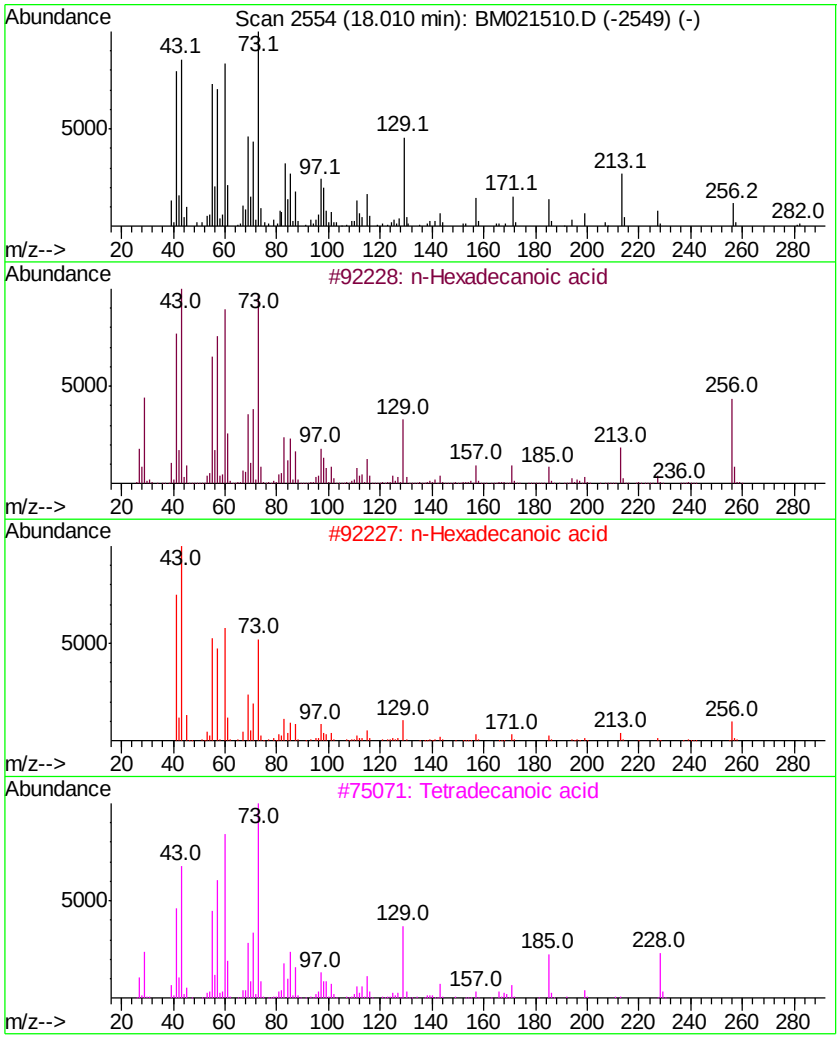
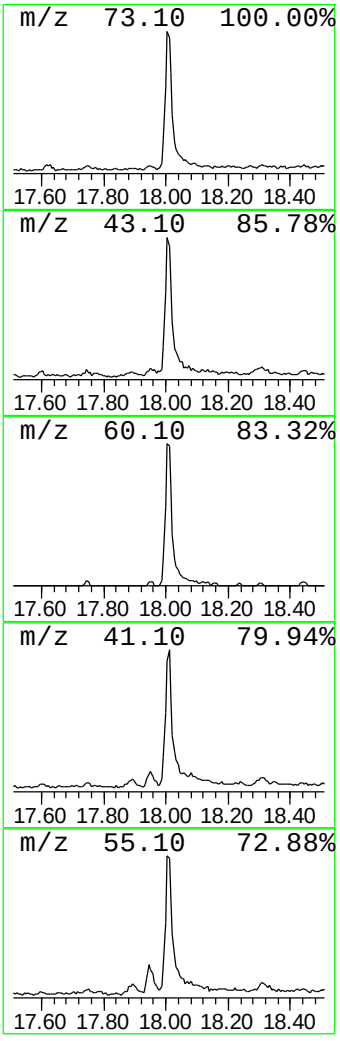
Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.01	5.30 ng/ul	455549	Phenanthrene-d10	17.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	70
4	n-Decanoic acid	172	C10H20O2	000334-48-5	70
5	n-Decanoic acid	172	C10H20O2	000334-48-5	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
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Acq On : 18 Jul 2019 04:22
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Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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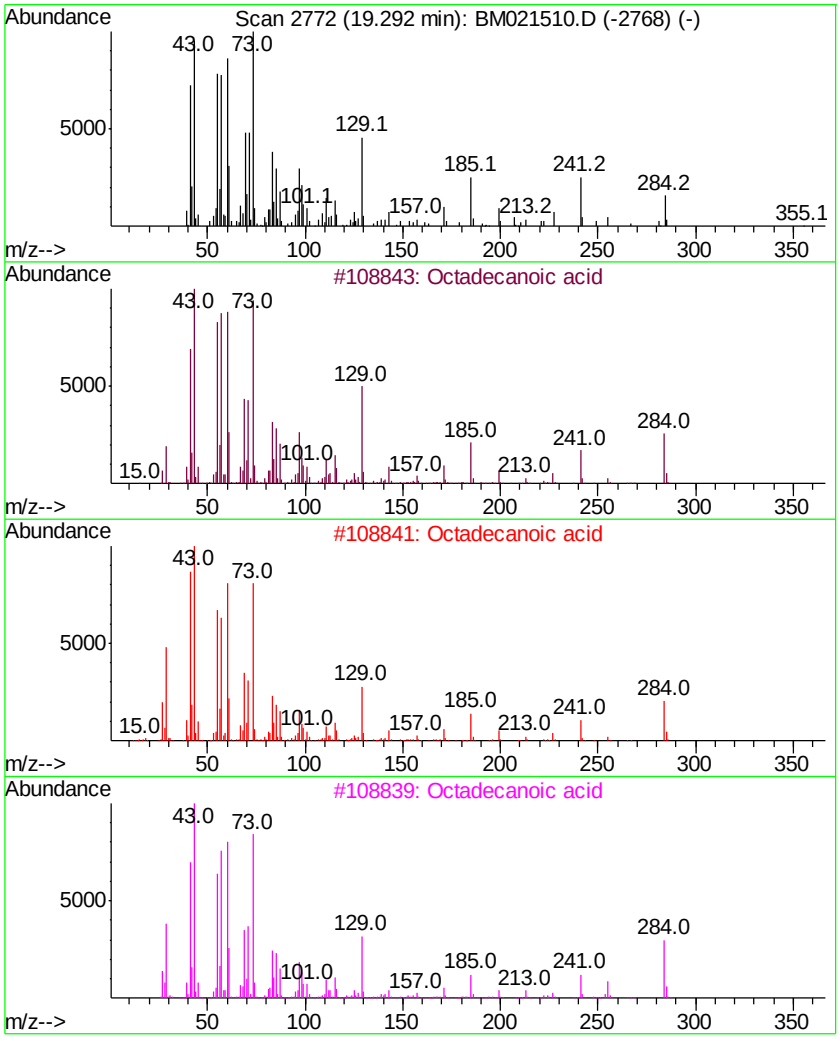
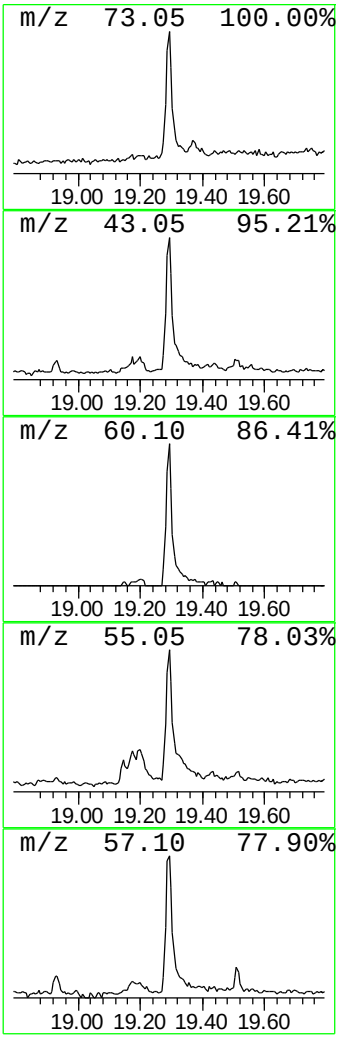
Peak Number 4 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.29	2.76 ng/ul	301439	Chrysene-d12	21.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	98
3	Octadecanoic acid	284	C18H36O2	000057-11-4	94
4	Octadecanoic acid	284	C18H36O2	000057-11-4	91
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	74





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Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
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Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

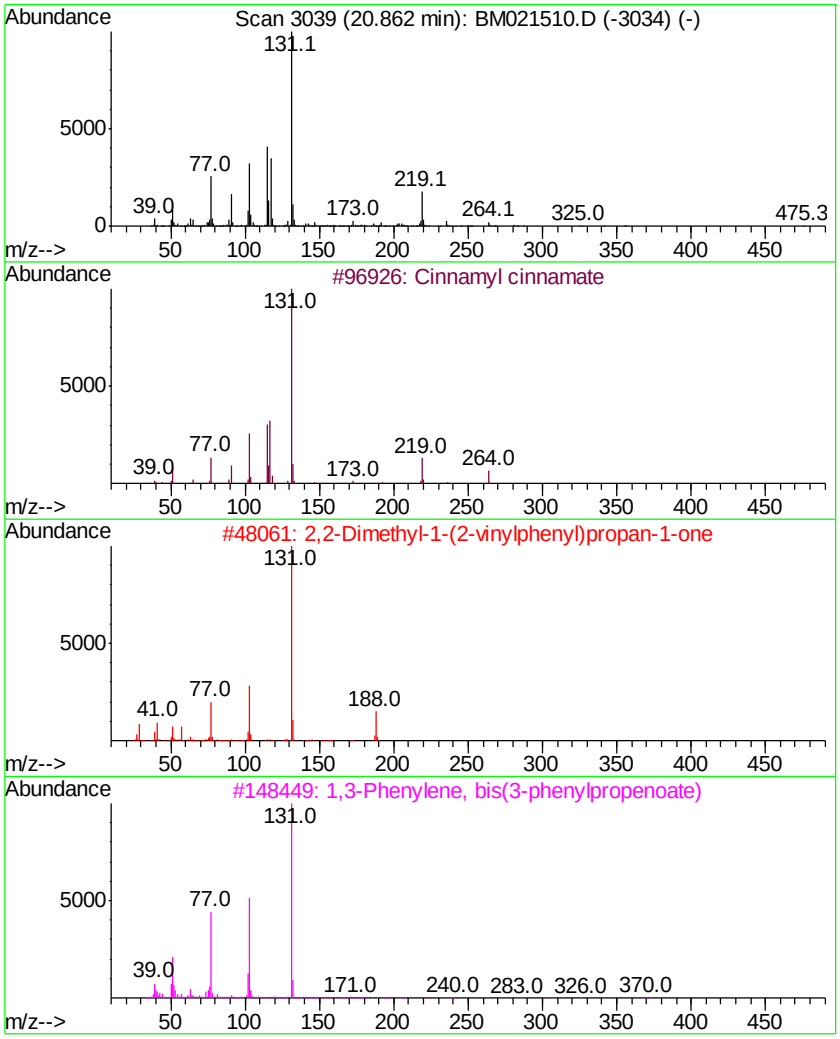
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

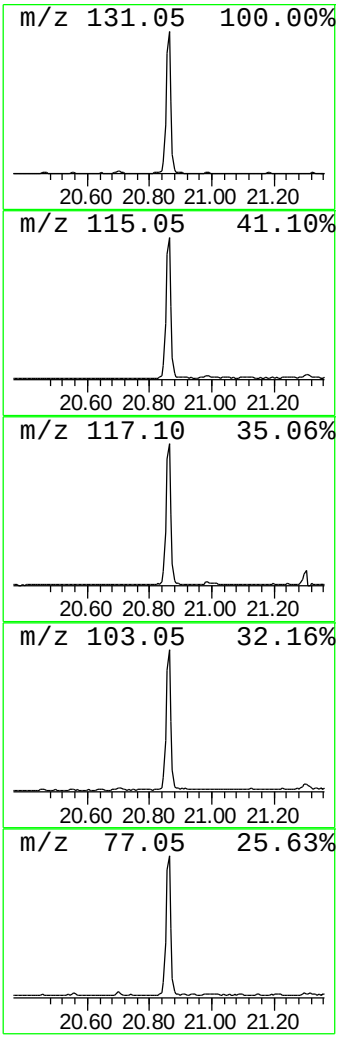
****

Peak Number 5 Cinnamyl cinnamate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	12.67 ng/ul	1382340	Chrysene-d12	21.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	86
2	2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	49
3	1,3-Phenylene, bis(3-phenylprope...	370	C24H18O4	1000259-62-4	47
4	1-Penten-3-one, 4-methyl-1-phenyl-	174	C12H14O	003160-32-5	47
5	1-Penten-3-one, 4,4-dimethyl-1-p...	188	C13H16O	000538-44-3	47





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
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Misc :
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Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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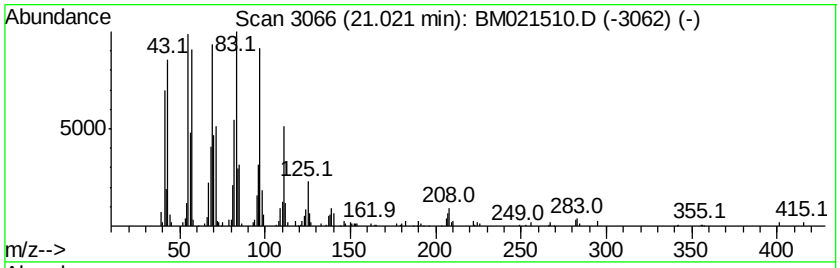
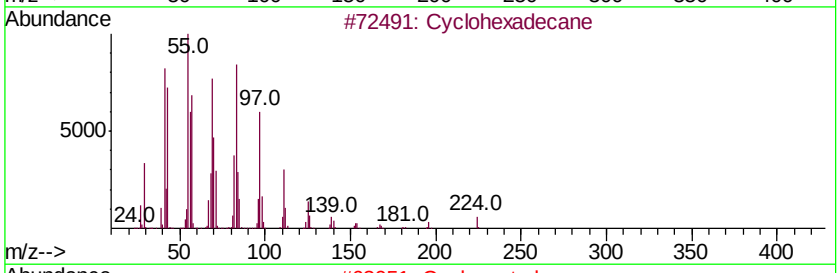
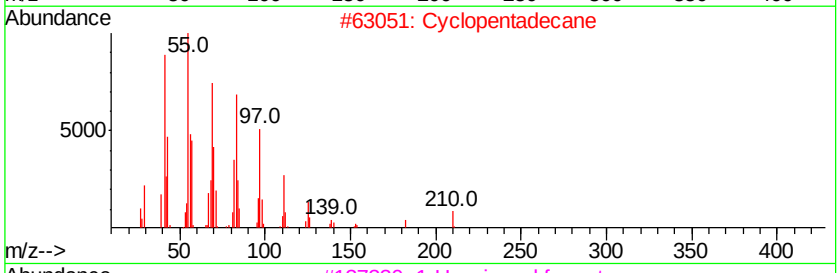
Peak Number 6 (DEL) Alkane: Cyclic21.02 Concentration Rank 8

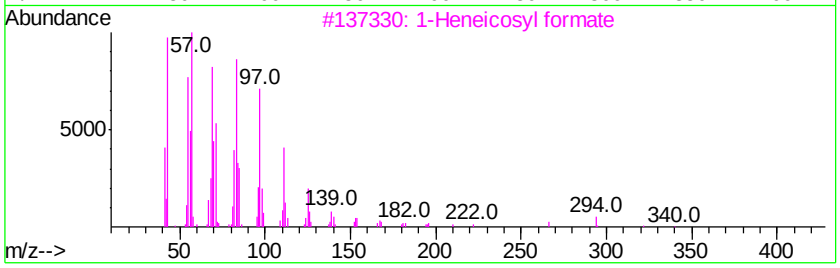
R.T.	EstConc	Area	Relative to ISTD	R.T.
21.02	4.94 ng/ul	538848	Chrysene-d12	21.30

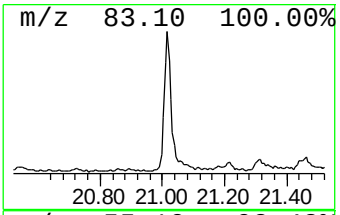
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	97
2	Cyclopentadecane	210	C15H30	000295-48-7	96
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
4	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	90
5	3-Eicosene, (E)-	280	C20H40	074685-33-9	90

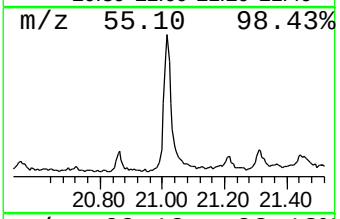


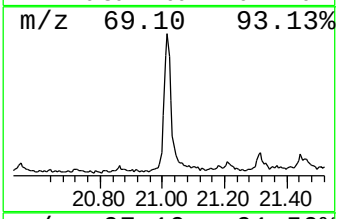


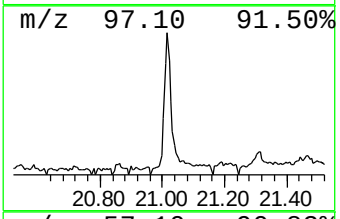


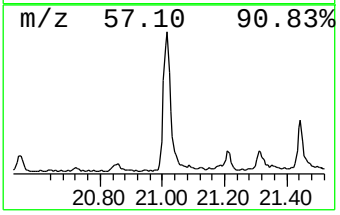












Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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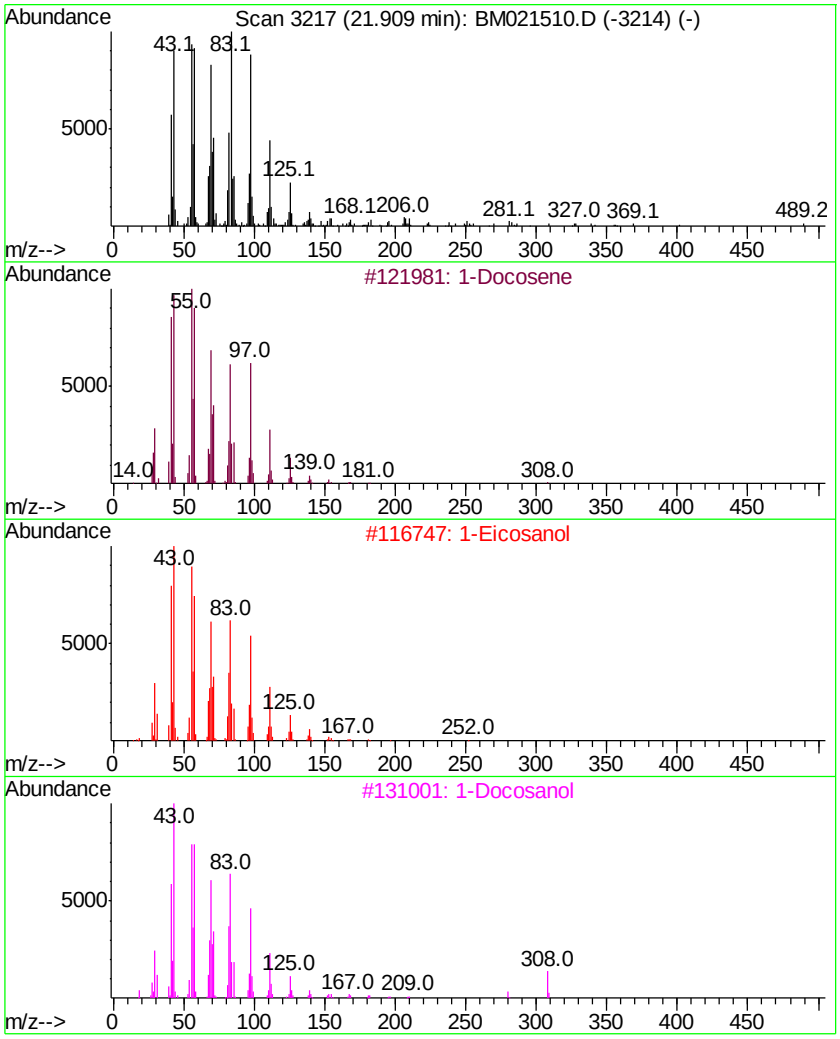
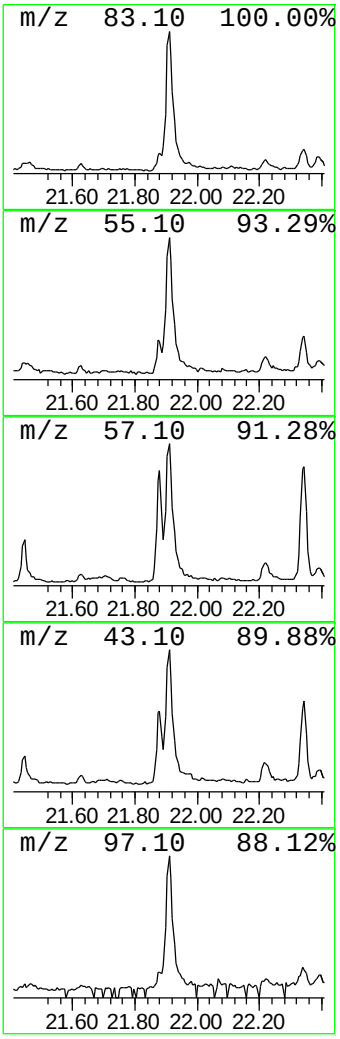
Peak Number 7 (DEL) Alkane: Cyclic21.91 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.91	5.58 ng/ul	609148	Chrysene-d12	21.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	86
2	1-Eicosanol	298	C20H42O	000629-96-9	81
3	1-Docosanol	326	C22H46O	000661-19-8	81
4	1-Nonadecene	266	C19H38	018435-45-5	81
5	1-Heptadecanol	256	C17H36O	001454-85-9	76





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Acq On : 18 Jul 2019 04:22
Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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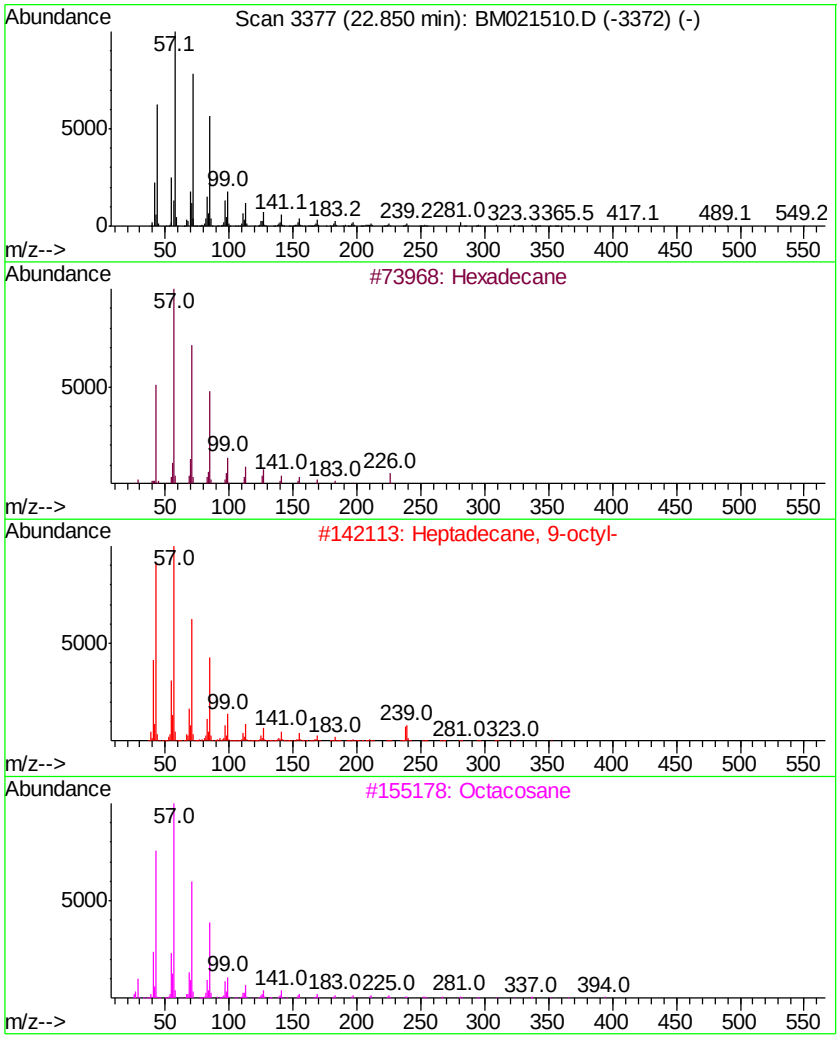
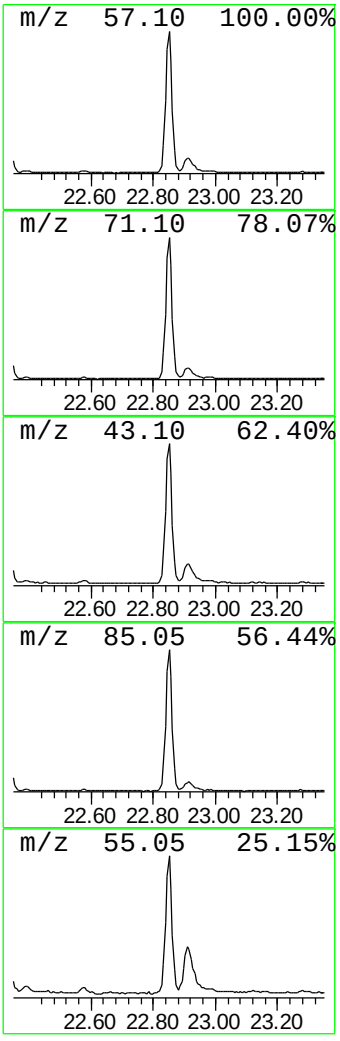
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.85	13.43 ng/ul	1527460	Perylene-d12	23.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Octacosane	394	C28H58	000630-02-4	91
4	Tetratetracontane	619	C44H90	007098-22-8	91
5	Heptacosane	380	C27H56	000593-49-7	91





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Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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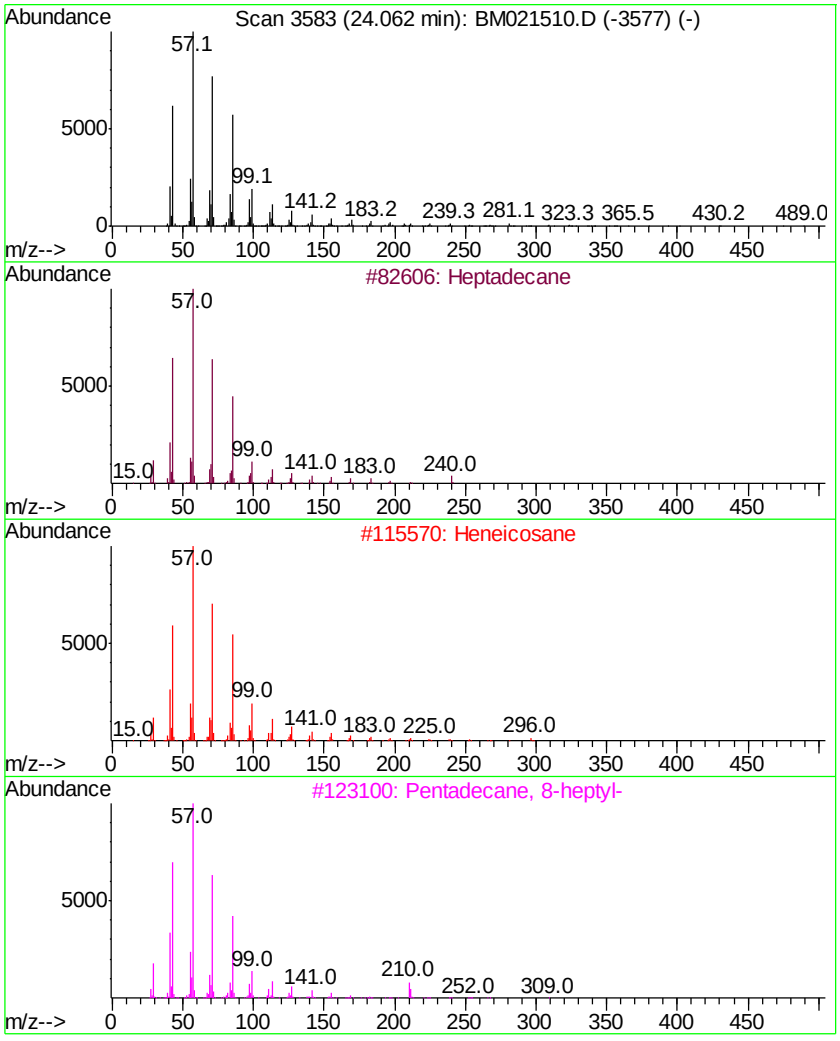
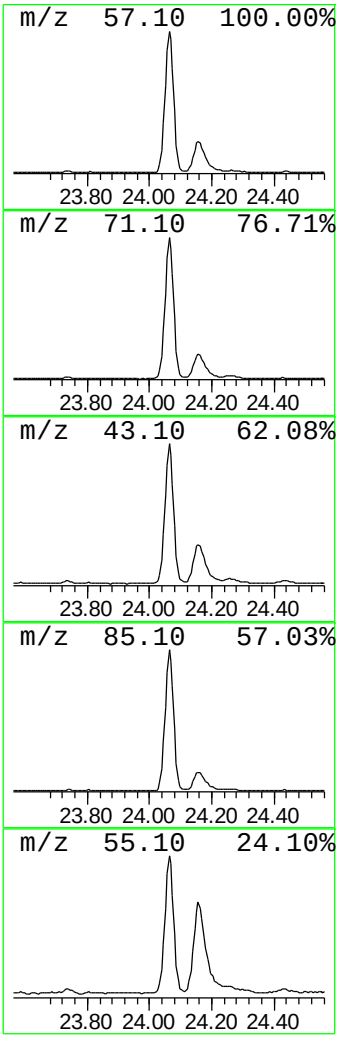
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.06	22.53 ng/ul	2562430	Perylene-d12	23.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	94
2	Heneicosane	296	C21H44	000629-94-7	91
3	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
4	Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
5	Nonacosane	408	C29H60	000630-03-5	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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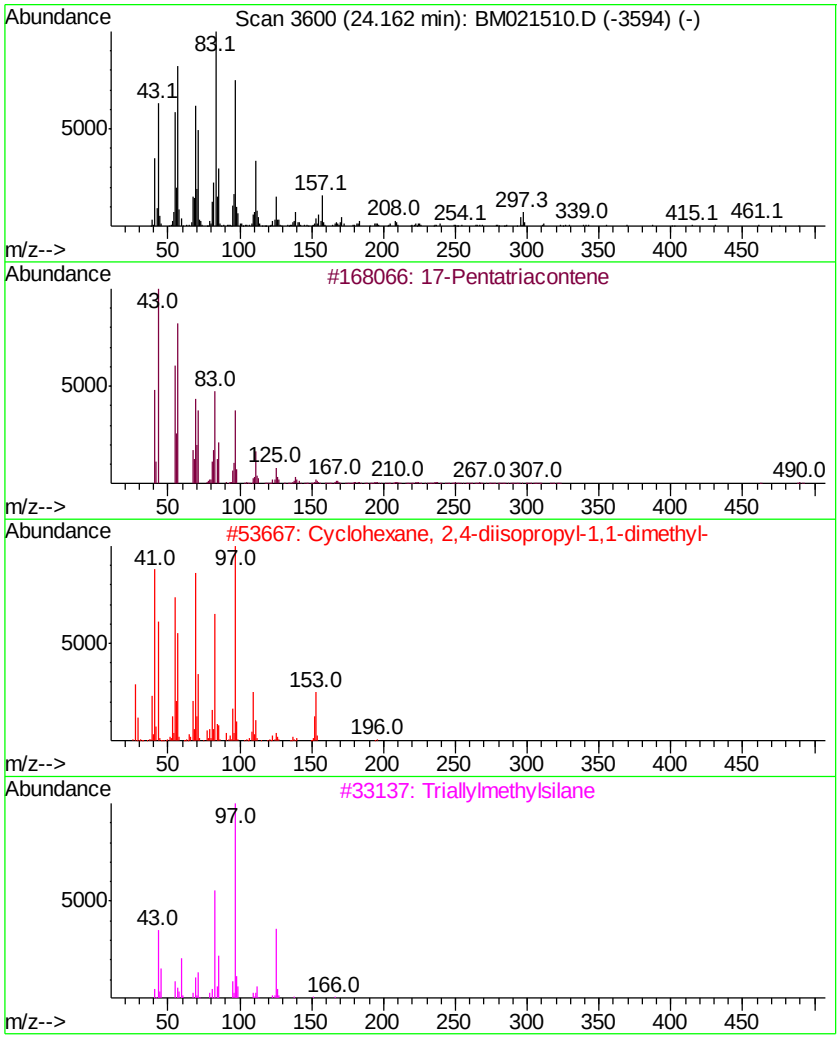
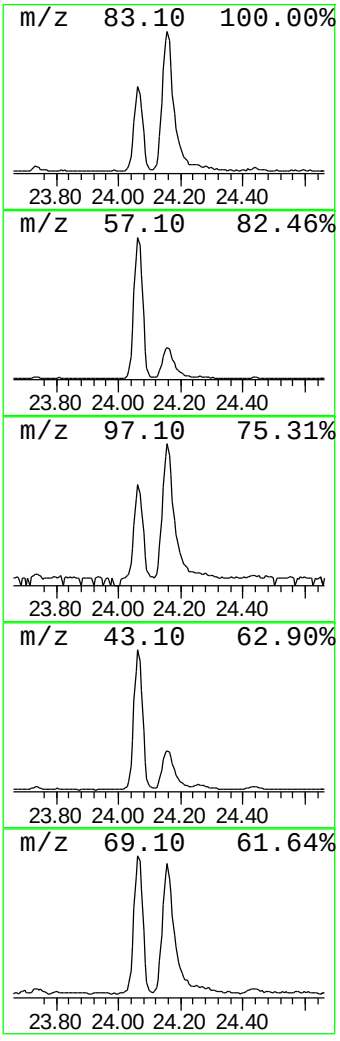
Peak Number 10 (DEL) Alkane: Cyclic24.16 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.16	12.81 ng/ul	1456710	Perylene-d12	23.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Pentatriacontene	491	C35H70	006971-40-0	53
2	Cyclohexane, 2,4-diisopropyl-1,1...	196	C14H28	1000149-60-5	49
3	Triallylmethylsilane	166	C10H18Si	001112-91-0	43
4	Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	43
5	2-Aziridinone, 1-tert-butyl-3-(1...	223	C14H25NO	026944-18-3	42





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071719\
Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION

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

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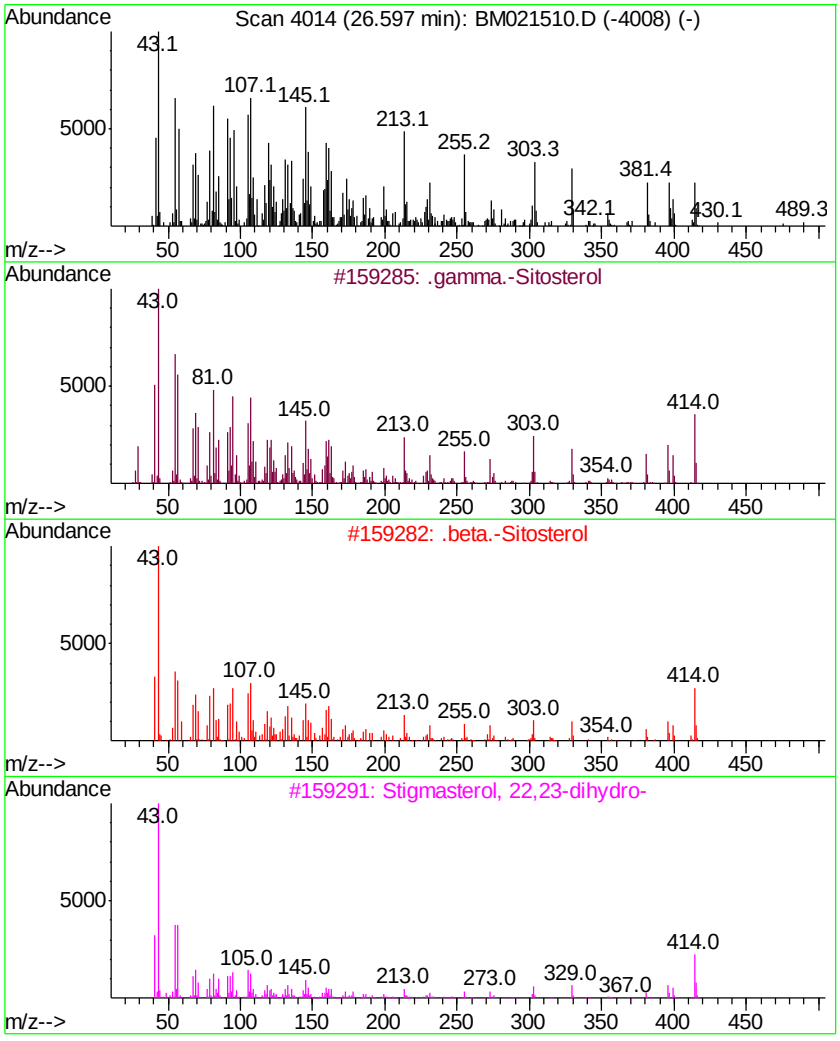
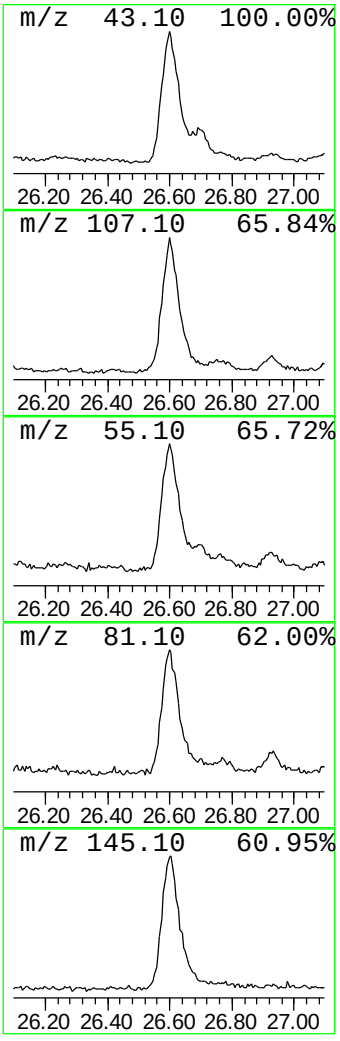
Peak Number 11 .gamma.-Sitosterol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.60	16.55 ng/ul	1882360	Perylene-d12	23.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	95
3	Stigmasterol, 22,23-dihydro-	414	C29H50O	1000214-20-7	86
4	.beta.-Sitosterol	414	C29H50O	000083-46-5	52
5	.beta.-Sitosterol	414	C29H50O	000083-46-5	38





Tentatively Identified Compound (LSC) summary

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Data File : BM021510.D
Acq On : 18 Jul 2019 04:22
Operator : HP/JU
Sample : K3760-08
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB8Y8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
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
Cyclotrisiloxane,...	4.37	2.7	ng/ul	108635	1	7.72	791605	20.0
Bicyclo[3.1.1]hep...	6.39	2.0	ng/ul	80392	1	7.72	791605	20.0
n-Hexadecanoic acid	18.01	5.3	ng/ul	455549	4	17.12	1719020	20.0
Octadecanoic acid	19.29	2.8	ng/ul	301439	5	21.30	2181970	20.0
Cinnamyl cinnamate	20.86	12.7	ng/ul	1382340	5	21.30	2181970	20.0
(DEL) Alkane: Cyc...	21.02	4.9	ng/ul	538848	5	21.30	2181970	20.0
(DEL) Alkane: Cyc...	21.91	5.6	ng/ul	609148	5	21.30	2181970	20.0
(DEL) Alkane: Str...	22.85	13.4	ng/ul	1527460	6	23.56	2274600	20.0
(DEL) Alkane: Str...	24.06	22.5	ng/ul	2562430	6	23.56	2274600	20.0
(DEL) Alkane: Cyc...	24.16	12.8	ng/ul	1456710	6	23.56	2274600	20.0
.gamma.-Sitosterol	26.60	16.6	ng/ul	1882360	6	23.56	2274600	20.0