LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
Data File : BM030823.D
Acq On : 06 Jul 2021 17:06
Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBK52

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\SFAM-EPA-BM070621.M
Title : SVOA CALIBRATION

Signal : TIC: BM030823.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.257	25	29	35	rBV	24128	34035	1.83%	0.107%
2	3.304	35	37	44	rVB2	8405	11365	0.61%	0.036%
3	3.410	53	55	59	rVB4	2796	3749	0.20%	0.012%
4	3.593	81	86	93	rBV3	10527	16450	0.89%	0.052%
5	3.669	96	99	122	rBV	227110	454649	24.51%	1.431%
6	3.928	140	143	147	rVB4	4917	6044	0.33%	0.019%
7	3.981	149	152	160	rVB2	7810	11678	0.63%	0.037%
8	4.316	205	209	218	rBV	18611	30252	1.63%	0.095%
9	4.634	259	263	267	rVB6	3402	4418	0.24%	0.014%
10	4.904	305	309	322	rBV	57677	92566	4.99%	0.291%
11	5.651	431	436	442	rBV	4291	7359	0.40%	0.023%
12	5.969	485	490	502	rVB	54533	84430	4.55%	0.266%
13	6.134	513	518	526	rBV	25661	39868	2.15%	0.125%
14	6.645	598	605	608	rBV6	2421	5460	0.29%	0.017%
15	6.763	619	625	630	rVB	18023	26938	1.45%	0.085%
16	6.934	647	654	662	rBV	384508	621322	33.49%	1.956%
17	7.104	676	683	694	rBV	288096	473321	25.52%	1.490%
18	7.298	710	716	727	rVB	399538	645857	34.82%	2.033%
19	7.410	733	735	740	rVB6	2171	2559	0.14%	0.008%
20	7.604	763	768	774	rBV	5112	8292	0.45%	0.026%
21	7.763	786	795	803	rVB2	264240	452917	24.42%	1.426%
22	7.886	811	816	822	rVB	9014	13752	0.74%	0.043%
23	7.986	828	833	843	rBV	34188	61666	3.32%	0.194%
24	8.069	843	847	852	rVB7	2571	4364	0.24%	0.014%
25	8.175	860	865	869	rBV4	2340	2971	0.16%	0.009%
26	8.239	871	876	881	rBV2	6906	11048	0.60%	0.035%
27	8.298	881	886	890	rBV	19368	31972	1.72%	0.101%
28	8.334	890	892	897	rVB3	7321	9144	0.49%	0.029%
29	8.469	901	915	925	rBV	434316	757468	40.83%	2.384%
30	8.586	928	935	942	rBV	405012	643557	34.69%	2.026%
31	8.669	947	949	955	rVB3	5832	7180	0.39%	0.023%
32	8.792	965	970	971	rBV4	2748	4379	0.24%	0.014%
33	8.833	971	977	985	rVV	213093	357769	19.29%	1.126%
34	8.922	985	992	996	rVV	354235	604527	32.59%	1.903%
35	8.963	996	999	1010	rVB	183239	290326	15.65%	0.914%
36	9.328	1056	1061	1065	rBV5	2418	3198	0.17%	0.010%

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Sample : M2905-01
Misc :
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Instrument : BNA_M
ClientSampleId : DBK52

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\SFAM-EPA-BM070621.M
Title : SVOA CALIBRATION

37	9.639	1107	1114	1124	rBV	285863	475778	25.65%	1.497%
38	9.722	1124	1128	1138	rVB4	14394	36580	1.97%	0.115%
39	10.175	1198	1205	1223	rBV	405254	733697	39.55%	2.309%
40	10.551	1259	1269	1273	rBV	338726	592221	31.93%	1.864%
41	10.598	1273	1277	1286	rVB	246203	402591	21.70%	1.267%
42	10.692	1286	1293	1310	rBV	167236	328613	17.71%	1.034%
43	11.257	1384	1389	1394	rVV2	10539	16415	0.88%	0.052%
44	12.074	1519	1528	1537	rBV	350852	626978	33.80%	1.973%
45	12.210	1545	1551	1561	rBV	252107	405562	21.86%	1.276%
46	12.433	1582	1589	1594	rBV2	6403	11474	0.62%	0.036%
47	12.580	1608	1614	1626	rVB	424604	673946	36.33%	2.121%
48	12.886	1662	1666	1671	rVB4	2561	3262	0.18%	0.010%
49	13.557	1776	1780	1781	rBV2	2091	2511	0.14%	0.008%
50	13.663	1794	1798	1800	rBV3	1451	1879	0.10%	0.006%
51	13.810	1816	1823	1842	rBV	695215	1021062	55.04%	3.214%
52	14.086	1861	1870	1880	rBV	844025	1237650	66.72%	3.895%
53	14.180	1885	1886	1890	rVB4	1981	2101	0.11%	0.007%
54	14.263	1897	1900	1904	rBV5	1496	2376	0.13%	0.007%
55	14.392	1916	1922	1928	rBV	506379	762480	41.10%	2.400%
56	14.457	1928	1933	1941	rVB	453410	659645	35.56%	2.076%
57	14.598	1951	1957	1983	rBV	180919	433054	23.35%	1.363%
58	14.827	1992	1996	1999	rBV3	4290	5873	0.32%	0.018%
59	15.386	2085	2091	2101	rBV	1013301	1517863	81.82%	4.777%
60	15.515	2107	2113	2128	rBV	263312	402276	21.69%	1.266%
61	15.651	2128	2136	2147	rVV	898165	1209565	65.21%	3.807%
62	16.168	2221	2224	2229	rBV5	3128	4325	0.23%	0.014%
63	16.233	2229	2235	2237	rBV4	2441	3438	0.19%	0.011%
64	16.262	2237	2240	2245	rVV7	1506	2520	0.14%	0.008%
65	16.392	2256	2262	2265	rBV5	1451	2902	0.16%	0.009%
66	16.615	2298	2300	2305	rVB5	1353	2047	0.11%	0.006%
67	16.927	2348	2353	2354	rBV3	1672	1870	0.10%	0.006%
68	16.956	2354	2358	2362	rVB	9052	12791	0.69%	0.040%
69	17.133	2382	2388	2392	rBV	583595	862276	46.48%	2.714%
70	17.174	2392	2395	2400	rVV	792083	1113834	60.04%	3.506%
71	17.233	2400	2405	2409	rVV	1048689	1584036	85.39%	4.986%
72	17.268	2409	2411	2422	rVB	358483	416458	22.45%	1.311%
73	17.545	2454	2458	2460	rVV2	3026	3825	0.21%	0.012%
74	17.609	2463	2469	2485	rVV	600917	878946	47.38%	2.766%
75	17.768	2492	2496	2499	rVV4	2038	3832	0.21%	0.012%
76	17.803	2499	2502	2506	rVB3	3060	4416	0.24%	0.014%

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Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBK52

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
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Title : SVOA CALIBRATION

77	18.027	2535	2540	2548	rBV	75085	126917	6.84%	0.399%
78	18.092	2548	2551	2556	rVB5	10323	17685	0.95%	0.056%
79	18.498	2615	2620	2625	rBV	10983	16711	0.90%	0.053%
80	18.556	2625	2630	2640	rVV	74234	118796	6.40%	0.374%
81	18.627	2640	2642	2648	rVB5	2165	2871	0.15%	0.009%
82	18.792	2666	2670	2675	rBV2	5728	6665	0.36%	0.021%
83	19.162	2729	2733	2736	rBV	9806	14605	0.79%	0.046%
84	19.309	2754	2758	2767	rBV5	29177	58016	3.13%	0.183%
85	19.409	2772	2775	2782	rVV3	6661	15166	0.82%	0.048%
86	19.521	2788	2794	2798	rBV	1197711	1855014	100.00%	5.839%
87	19.550	2798	2799	2806	rVB	492274	445771	24.03%	1.403%
88	19.756	2831	2834	2838	rBV2	15556	18623	1.00%	0.059%
89	21.039	3048	3052	3057	rBV	46323	66199	3.57%	0.208%
90	21.092	3057	3061	3068	rBV	119959	154615	8.33%	0.487%
91	21.309	3093	3098	3101	rBV	632242	805675	43.43%	2.536%
92	21.344	3101	3104	3110	rVB	760748	938727	50.60%	2.955%
93	22.880	3359	3365	3369	rBV	450912	784571	42.29%	2.469%
94	22.927	3369	3373	3383	rVB	329463	561272	30.26%	1.767%
95	23.415	3450	3456	3471	rVB2	804973	1815080	97.85%	5.713%
96	23.556	3475	3480	3489	rVB2	406724	778584	41.97%	2.451%
97	24.032	3557	3561	3566	rBV	106094	191354	10.32%	0.602%
98	26.532	3979	3986	4001	rVB	214646	678964	36.60%	2.137%

Sum of corrected areas: 31771699

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Data File : BM030823.D
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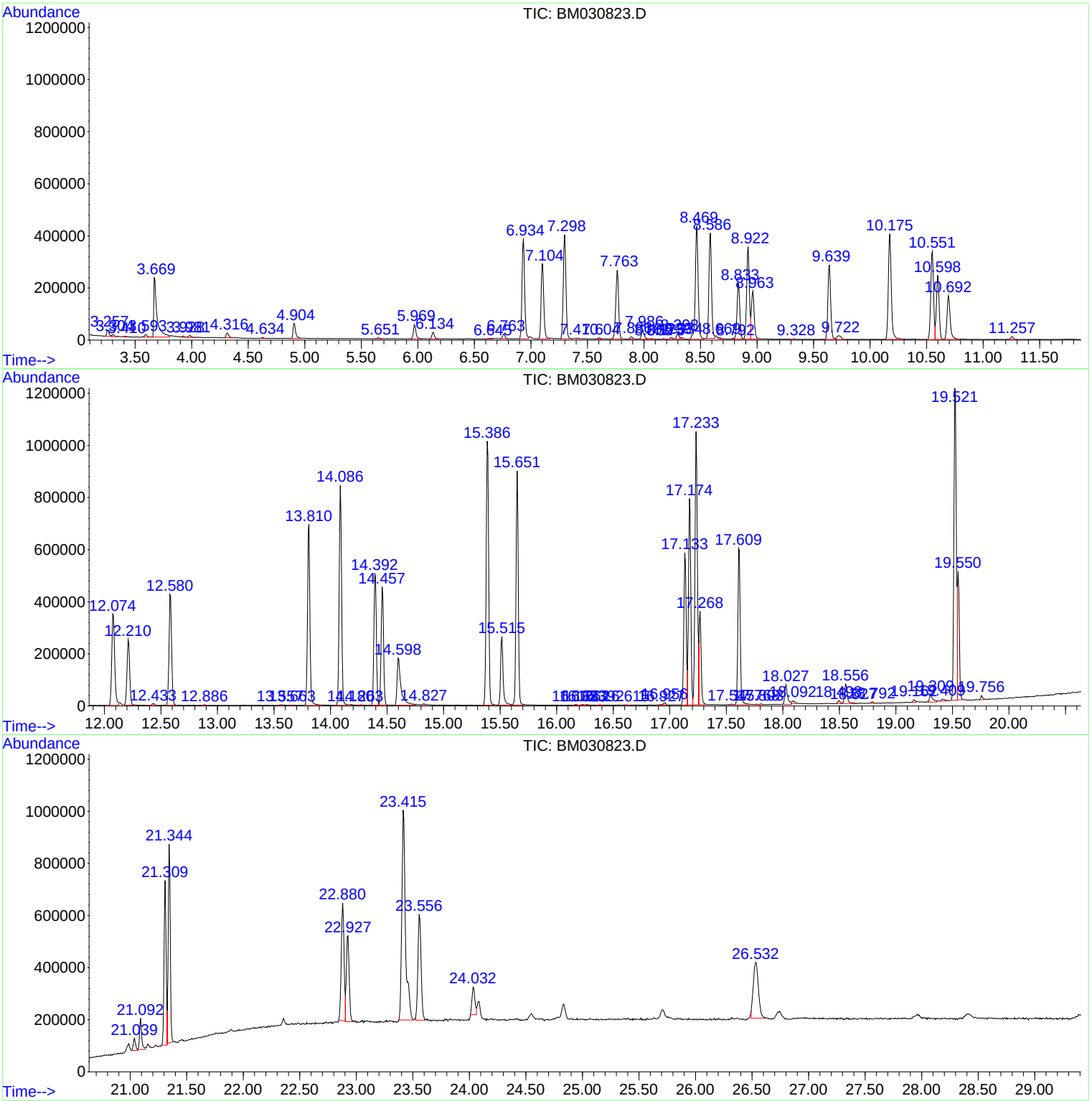
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBK52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
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Instrument :
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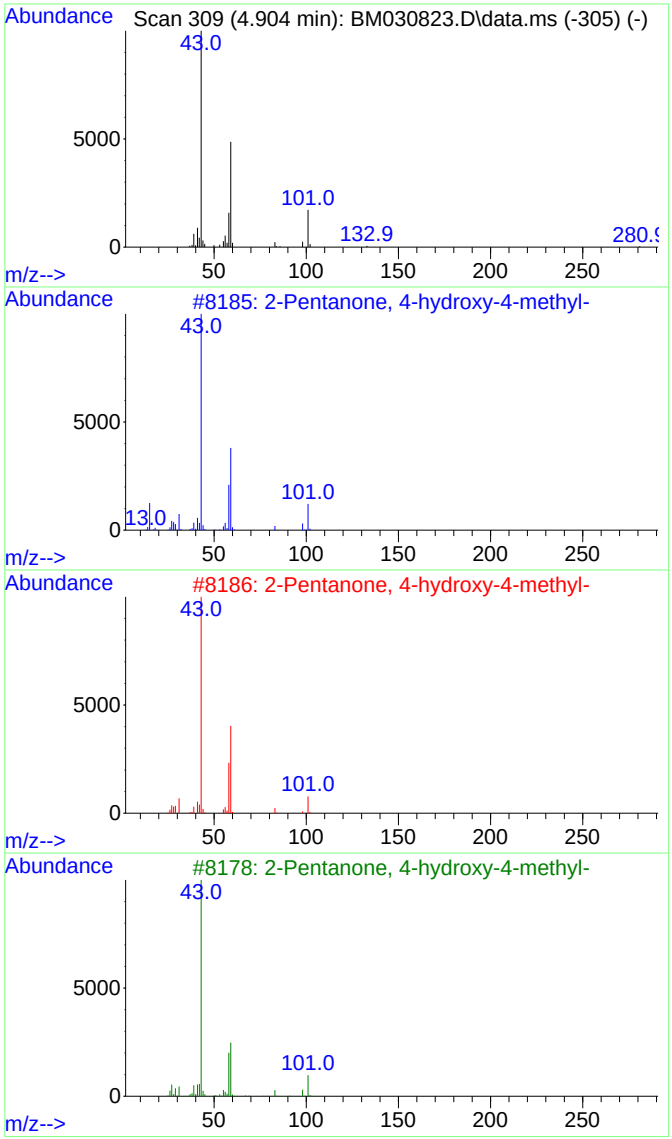
Quant Title : SVOA CALIBRATION

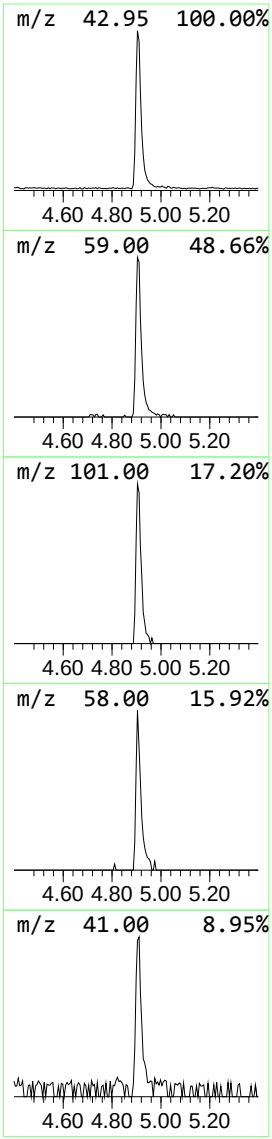
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.900	4.09 ng/ul	92566	1,4-Dichlorobenzene-d4	7.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
Data File : BM030823.D
Acq On : 06 Jul 2021 17:06
Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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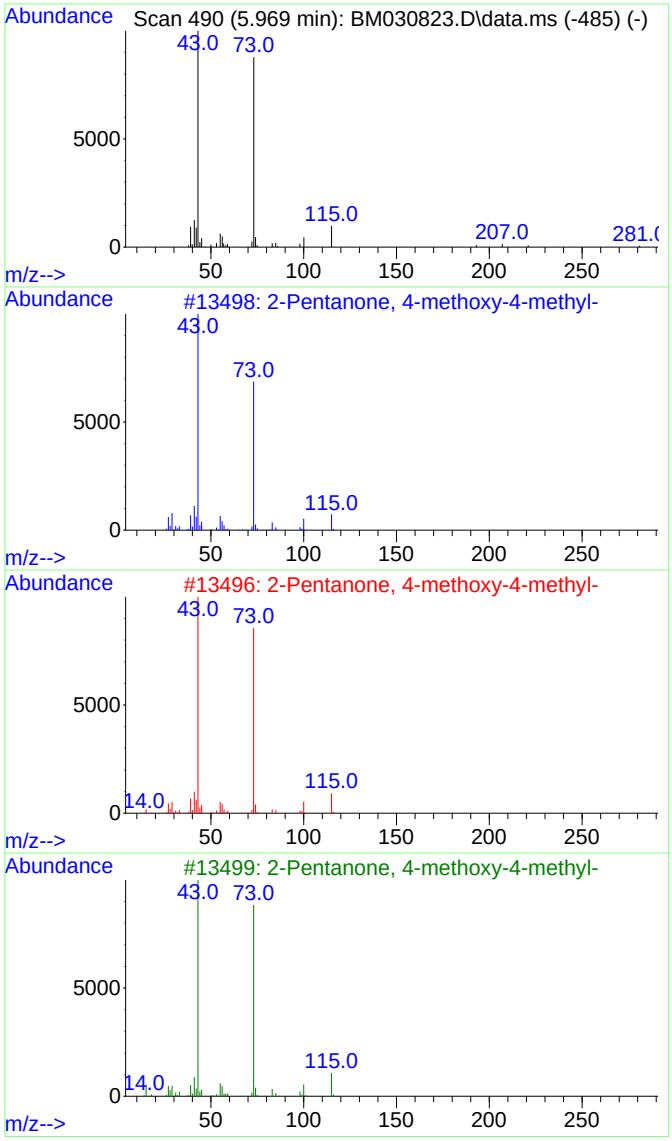
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

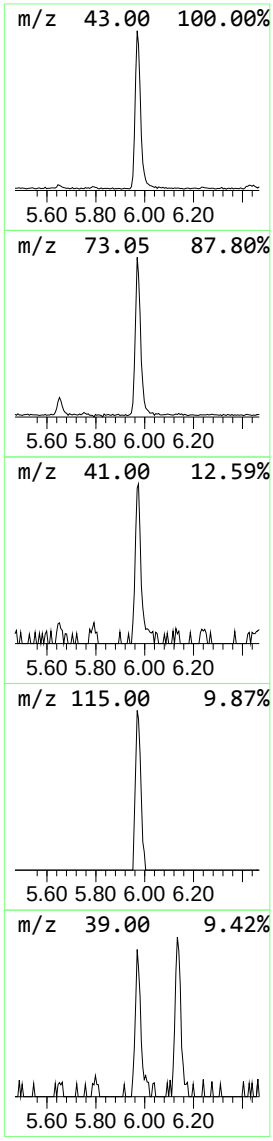
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.970	3.73 ng/ul	84430	1,4-Dichlorobenzene-d4	7.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78
2	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78
3	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	72
4	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	64
5	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	37





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BNA_M
ClientSampleId :
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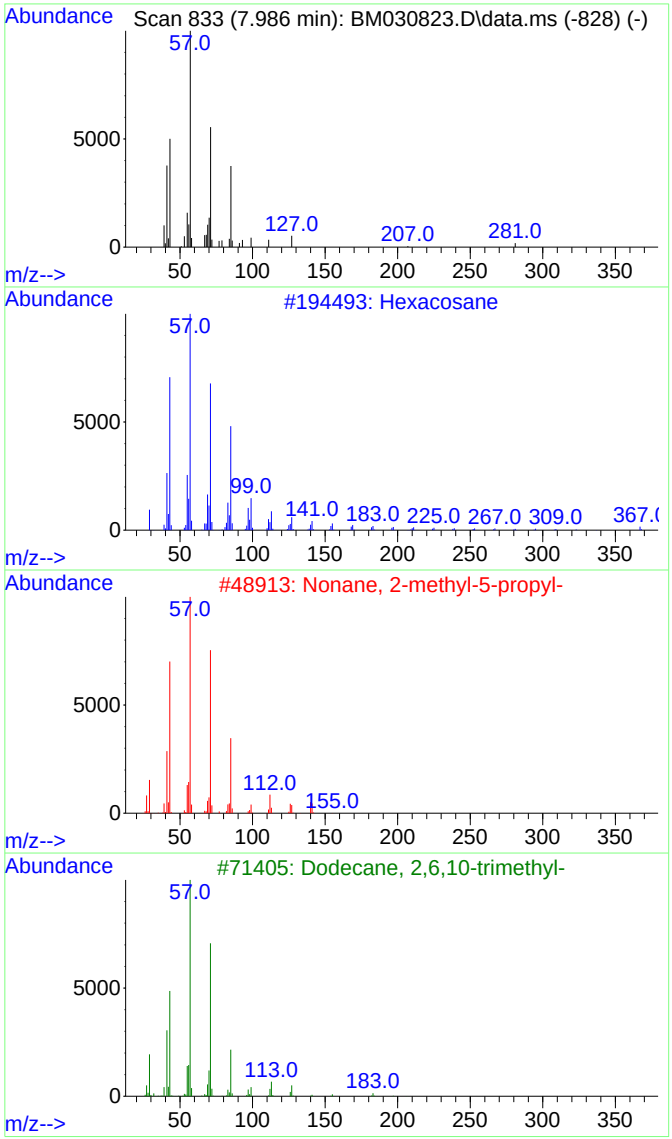
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

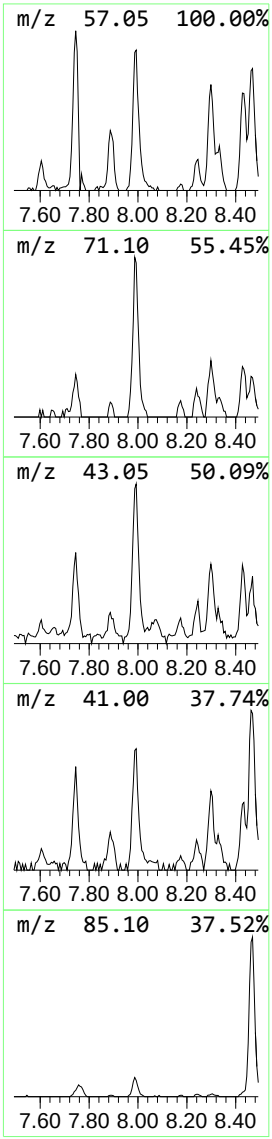
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.990	2.72 ng/ul	61666	1,4-Dichlorobenzene-d4	7.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	72
2	Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	72
3	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
4	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
5	Decane, 5-propyl-	184	C13H28	017312-62-8	64





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Quant Title : SVOA CALIBRATION

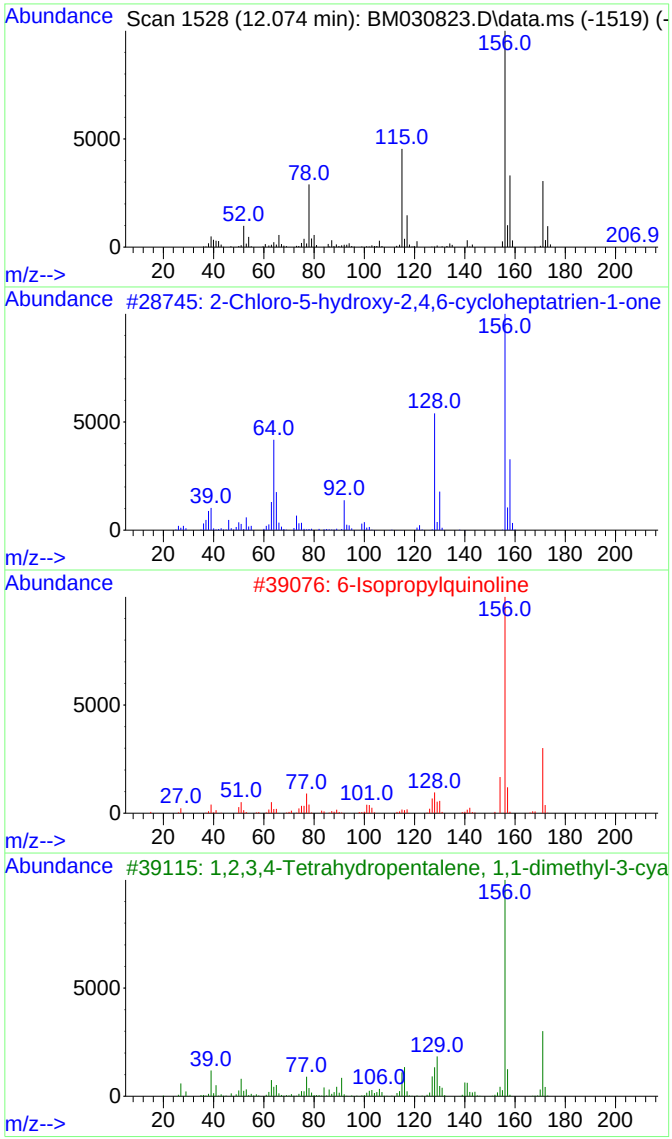
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

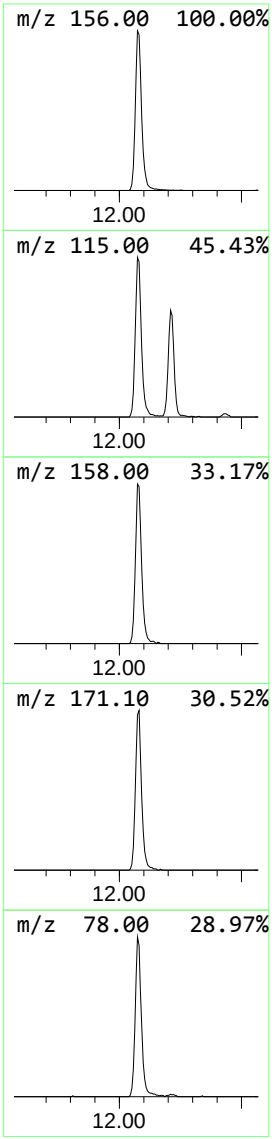
****

Peak Number 4 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.070	21.17 ng/ul	626978	Naphthalene-d8	10.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2	6-Isopropylquinoline	171	C12H13N	000135-79-5	38
3	1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	37
4	Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32
5	Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
Data File : BM030823.D
Acq On : 06 Jul 2021 17:06
Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBK52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

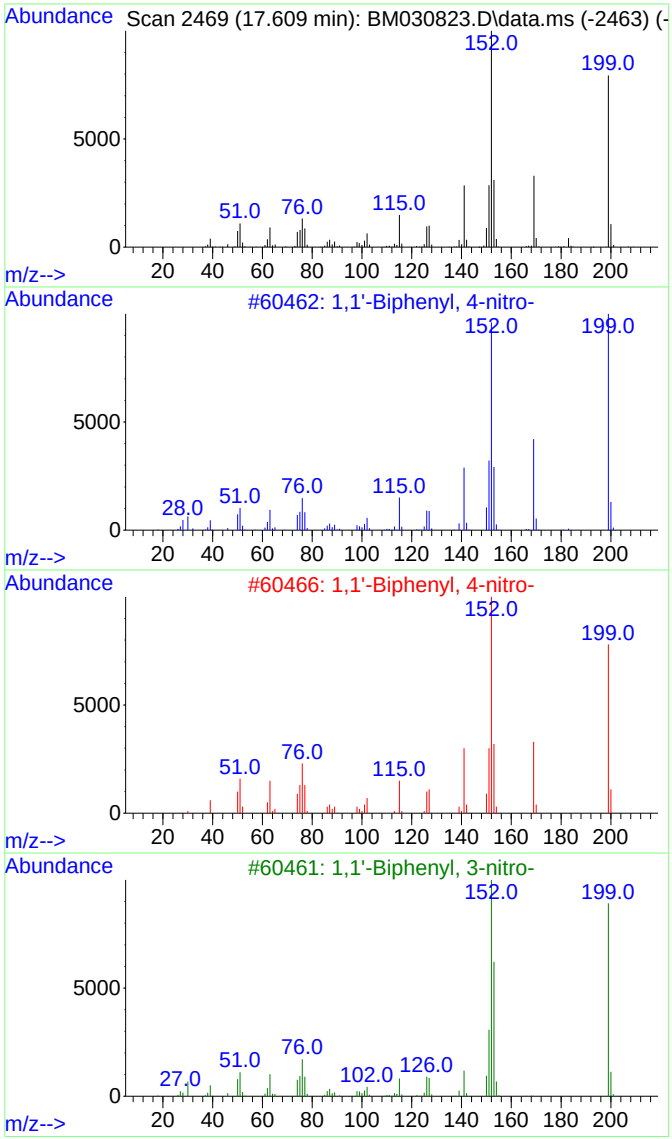
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

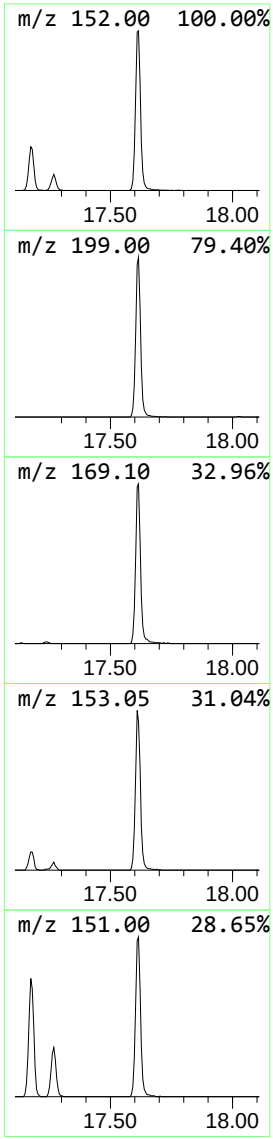
****

Peak Number 5 1,1'-Biphenyl, 4-nitro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.610	20.39 ng/ul	878946	Phenanthrene-d10	17.133

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	99
2	1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
3	1,1'-Biphenyl, 3-nitro-	199	C12H9NO2	002113-58-8	90
4	Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	87
5	Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	86





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Data File : BM030823.D
Acq On : 06 Jul 2021 17:06
Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBK52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

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

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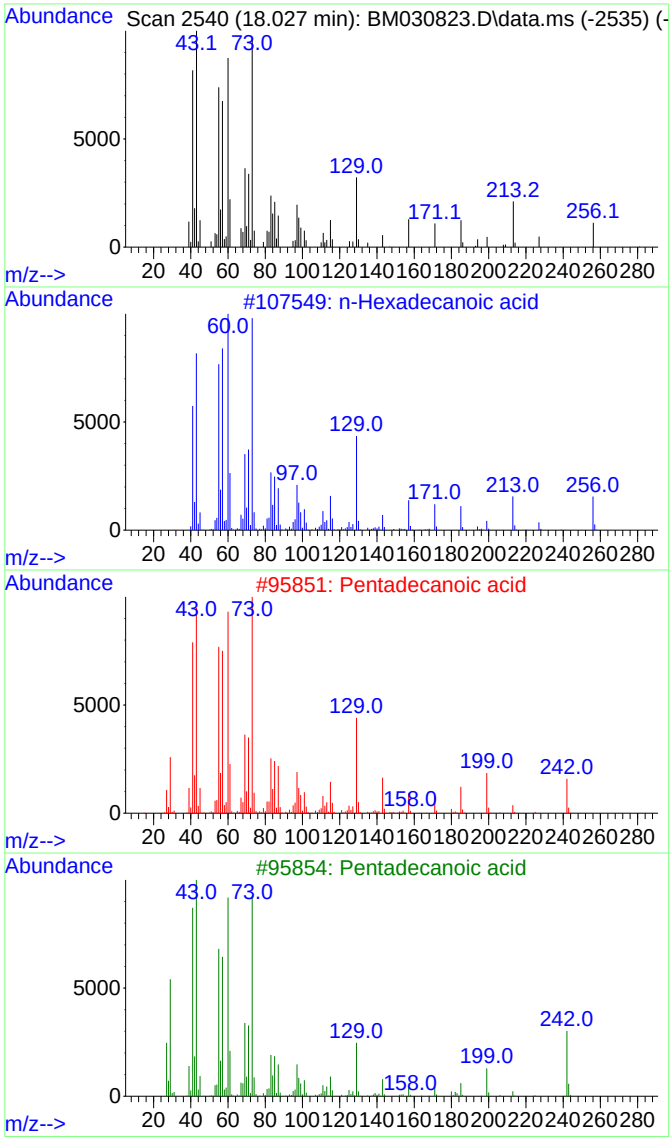
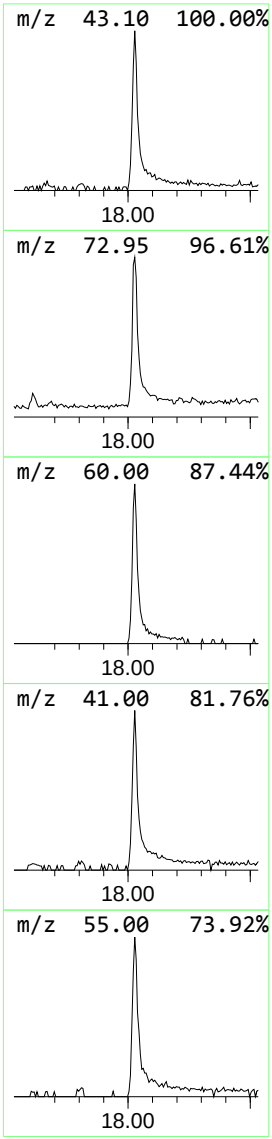
Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.030	2.94 ng/ul	126917	Phenanthrene-d10	17.133

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	81
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	64





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Data File : BM030823.D
Acq On : 06 Jul 2021 17:06
Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBK52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

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

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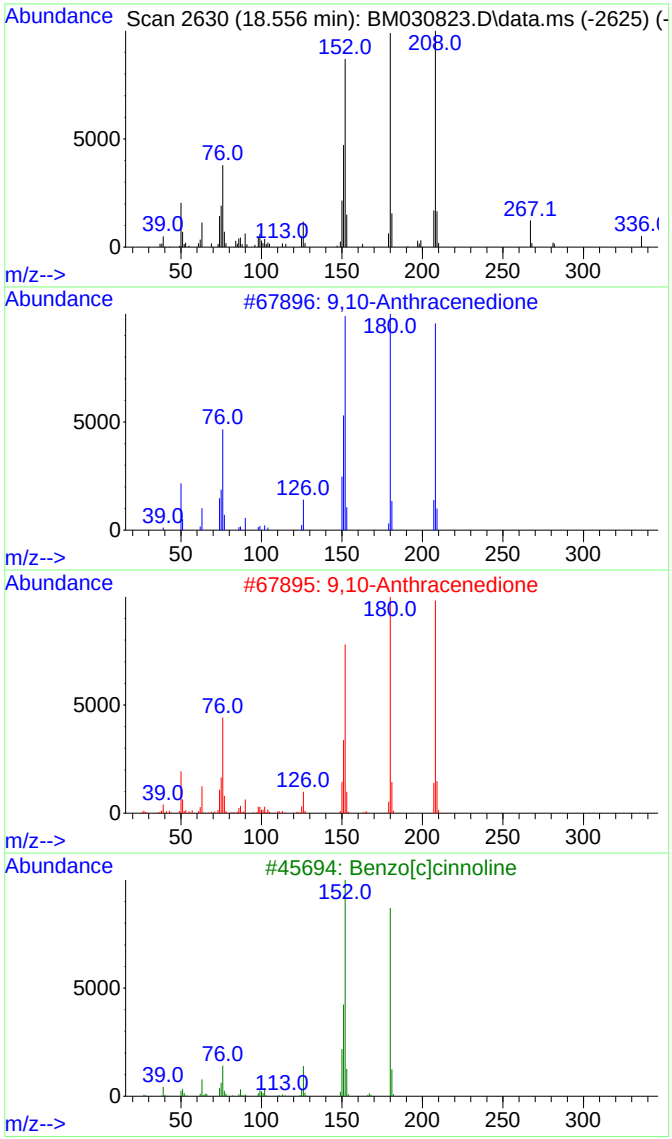
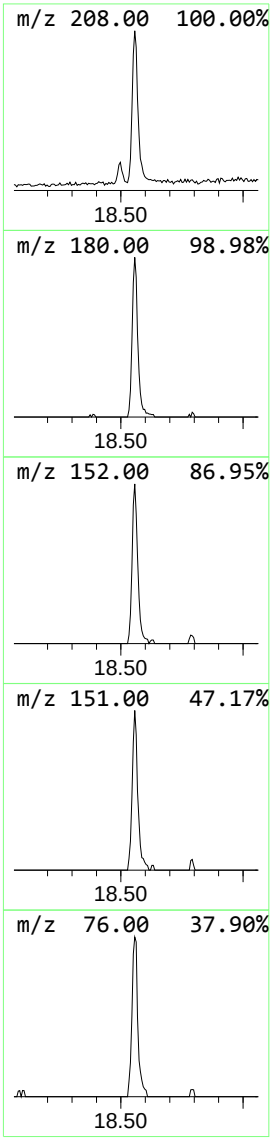
Peak Number 7 9,10-Anthracenedione Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.560	2.76 ng/ul	118796	Phenanthrene-d10	17.133

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
3	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	96
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	83





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Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBK52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

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

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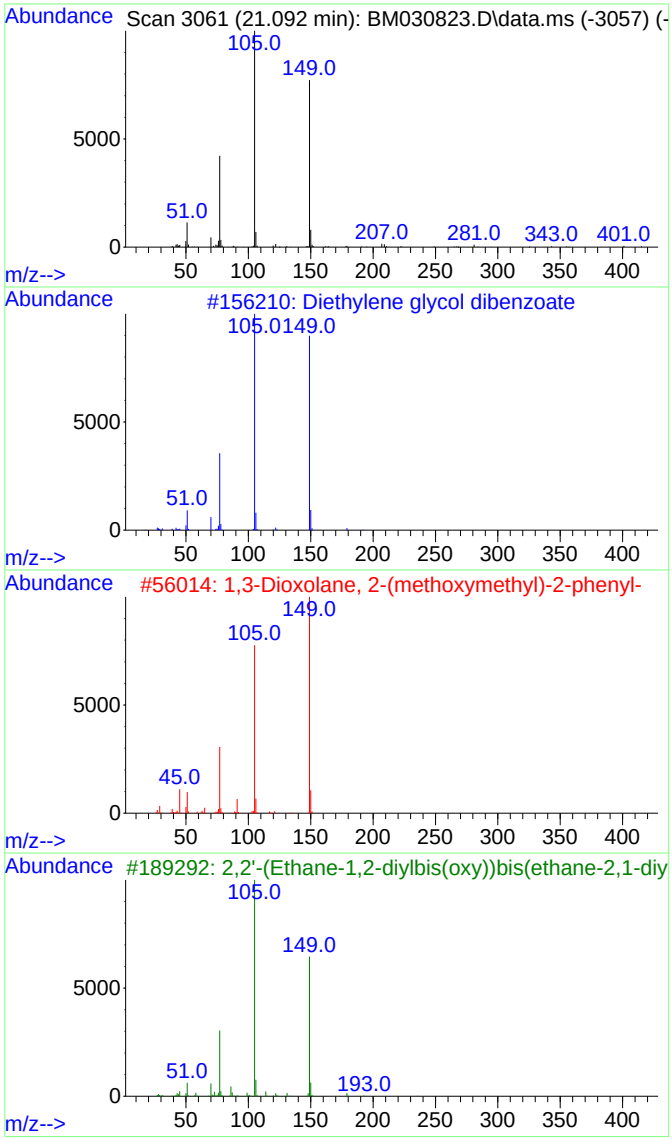
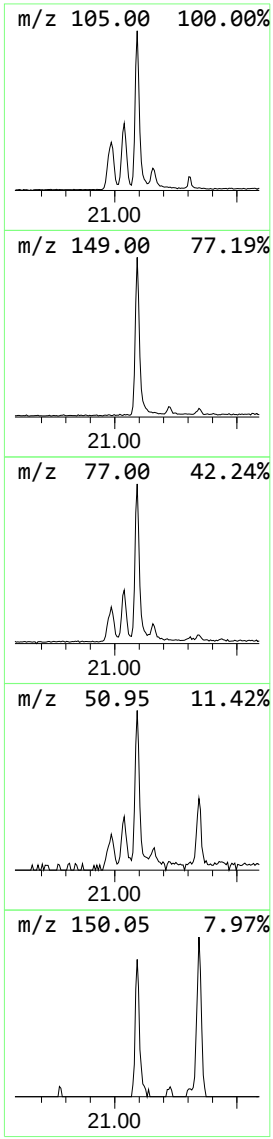
Peak Number 8 Diethylene glycol dibenzoate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.090	3.84 ng/ul	154615	Chrysene-d12	21.309

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	83
3	2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	83
4	2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	52
5	2-Bromoisobutyrophenone	226	C10H11BrO	010409-54-8	35





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Data File : BM030823.D
Acq On : 06 Jul 2021 17:06
Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBK52

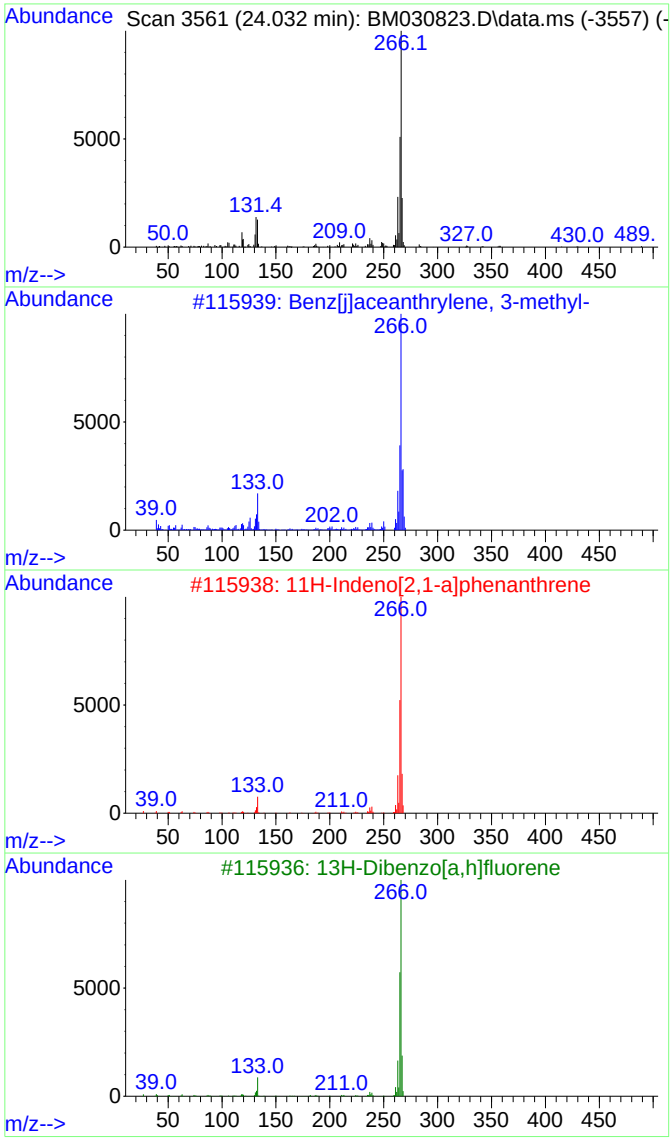
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.M
Quant Title : SVOA CALIBRATION

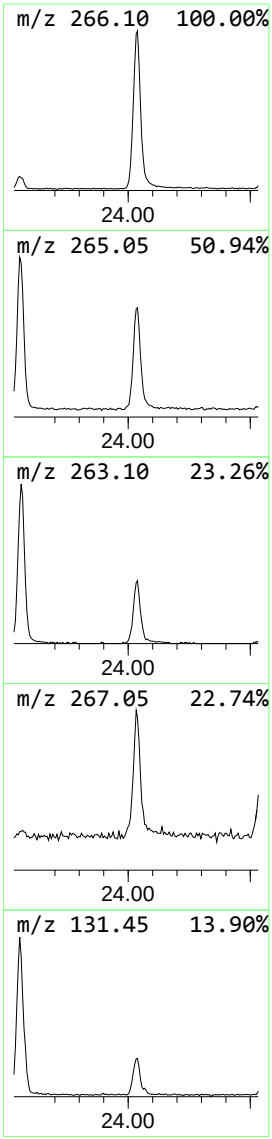
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benz[j]aceanthrylene, 3-met... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.030	4.92 ng/ul	191354	Perylene-d12	23.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[j]aceanthrylene, 3-methyl-	266	C21H14	003343-10-0	92
2	11H-Indeno[2,1-a]phenanthrene	266	C21H14	000220-97-3	90
3	13H-Dibenzo[a,h]fluorene	266	C21H14	000239-85-0	81
4	8H-Indeno[2,1-b]phenanthrene	266	C21H14	000241-28-1	64
5	Perylene, 3-methyl-	266	C21H14	024471-47-4	58





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070621\
Data File : BM030823.D
Acq On : 06 Jul 2021 17:06
Operator : CG/JU
Sample : M2905-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DBK52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM070621.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
2-Pentanone, 4-...	4.900	4.1	ng/ul	92566	1	7.763	452917	20.0
2-Pentanone, 4-...	5.970	3.7	ng/ul	84430	1	7.763	452917	20.0
(DEL) Alkane: S...	7.990	2.7	ng/ul	61666	1	7.763	452917	20.0
unknown-01	12.070	21.2	ng/ul	626978	2	10.551	592221	20.0
1,1'-Biphenyl, ...	17.610	20.4	ng/ul	878946	4	17.133	862276	20.0
n-Hexadecanoic ...	18.030	2.9	ng/ul	126917	4	17.133	862276	20.0
9,10-Anthracene...	18.560	2.8	ng/ul	118796	4	17.133	862276	20.0
Diethylene glyc...	21.090	3.8	ng/ul	154615	5	21.309	805675	20.0
Benz[j]aceanthr...	24.030	4.9	ng/ul	191354	6	23.556	778584	20.0