LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBYF1

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_G\Methods\SFAM-EPA-BG122422.M
Title : SVOA CALIBRATION

Signal : TIC: BG056103.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.370	9	14	28	rVB	594242	889907	94.76%	7.340%
2	3.652	57	62	65	rBV3	3684	5148	0.55%	0.042%
3	3.905	102	105	107	rBV3	1098	1776	0.19%	0.015%
4	4.093	135	137	141	rBV2	8134	9543	1.02%	0.079%
5	4.199	151	155	160	rVB3	4511	6510	0.69%	0.054%
6	4.328	173	177	182	rBV2	4345	6704	0.71%	0.055%
7	4.440	193	196	201	rVB2	2372	2562	0.27%	0.021%
8	4.774	249	253	254	rVB2	1401	1535	0.16%	0.013%
9	5.015	289	294	296	rBV2	1432	2009	0.21%	0.017%
10	5.039	296	298	304	rVB2	4385	5445	0.58%	0.045%
11	5.380	351	356	364	rVB3	14411	23896	2.54%	0.197%
12	5.997	456	461	464	rBV3	1130	1691	0.18%	0.014%
13	7.471	704	712	723	rBV	77334	138764	14.78%	1.145%
14	7.653	737	743	752	rVB	53259	89011	9.48%	0.734%
15	7.871	773	780	786	rBV	72961	122479	13.04%	1.010%
16	8.347	853	861	868	rBB	111802	194443	20.70%	1.604%
17	8.564	894	898	903	rBV2	1534	3134	0.33%	0.026%
18	9.034	971	978	985	rBV2	99390	169402	18.04%	1.397%
19	9.175	996	1002	1009	rBV2	8492	15163	1.61%	0.125%
20	9.516	1053	1060	1068	rVB2	63035	113695	12.11%	0.938%
21	9.898	1121	1125	1130	rVB2	2831	3867	0.41%	0.032%
22	10.245	1177	1184	1191	rBV	53092	92651	9.87%	0.764%
23	10.785	1268	1276	1283	rVB	120798	212741	22.65%	1.755%
24	11.179	1334	1343	1351	rBB	157216	295220	31.43%	2.435%
25	11.296	1358	1363	1372	rBV2	35534	70578	7.51%	0.582%
26	11.696	1426	1431	1432	rBV2	1289	1567	0.17%	0.013%
27	12.659	1590	1595	1600	rBV2	14980	22960	2.44%	0.189%
28	12.736	1604	1608	1612	rVB2	2089	2589	0.28%	0.021%
29	13.470	1729	1733	1739	rBV2	1673	2213	0.24%	0.018%
30	13.740	1777	1779	1781	rBV2	1856	1959	0.21%	0.016%
31	13.870	1797	1801	1804	rVB2	1489	1737	0.18%	0.014%
32	13.934	1804	1812	1814	rBV2	1120	2246	0.24%	0.019%
33	13.975	1817	1819	1824	rVB2	1900	2270	0.24%	0.019%
34	14.116	1838	1843	1847	rBV3	4402	7373	0.79%	0.061%
35	14.234	1859	1863	1867	rBV2	13309	16464	1.75%	0.136%
36	14.281	1867	1871	1876	rVV4	6320	8811	0.94%	0.073%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBYF1

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_G\Methods\SFAM-EPA-BG122422.M
Title : SVOA CALIBRATION

37	14.340	1876	1881	1888	rVB	245804	362805	38.63%	2.993%
38	14.475	1899	1904	1909	rBV3	5220	9040	0.96%	0.075%
39	14.651	1927	1934	1943	rVB	276554	448068	47.71%	3.696%
40	14.898	1972	1976	1977	rBV2	2707	2369	0.25%	0.020%
41	14.951	1979	1985	1993	rVB	352054	553608	58.95%	4.566%
42	15.109	2007	2012	2022	rBV2	87099	140598	14.97%	1.160%
43	15.198	2022	2027	2035	rVB3	4796	9197	0.98%	0.076%
44	15.268	2035	2039	2044	rBV4	3962	5838	0.62%	0.048%
45	15.321	2044	2048	2049	rVV	2767	2195	0.23%	0.018%
46	15.344	2049	2052	2056	rVB	2295	2258	0.24%	0.019%
47	15.938	2146	2153	2163	rVB	418518	638949	68.03%	5.270%
48	16.038	2163	2170	2176	rBV	83519	125783	13.39%	1.038%
49	16.120	2182	2184	2187	rVB	1690	1529	0.16%	0.013%
50	16.179	2190	2194	2198	rVV	2299	3734	0.40%	0.031%
51	16.226	2198	2202	2207	rVV4	15455	22474	2.39%	0.185%
52	16.290	2207	2213	2220	rVB	203145	287269	30.59%	2.370%
53	16.384	2227	2229	2233	rBV	1601	1560	0.17%	0.013%
54	16.426	2233	2236	2242	rVB3	4800	5815	0.62%	0.048%
55	16.478	2242	2245	2246	rBV3	1702	1443	0.15%	0.012%
56	16.561	2256	2259	2260	rBV	1573	1390	0.15%	0.011%
57	16.590	2262	2264	2268	rVB	1752	2459	0.26%	0.020%
58	16.708	2282	2284	2287	rBV2	1811	1459	0.16%	0.012%
59	16.860	2305	2310	2314	rVB3	5500	8859	0.94%	0.073%
60	17.042	2337	2341	2347	rVB2	2363	3908	0.42%	0.032%
61	17.289	2378	2383	2385	rVB	1531	1729	0.18%	0.014%
62	17.342	2387	2392	2394	rBV	1310	1646	0.18%	0.014%
63	17.471	2413	2414	2418	rBV	1713	1390	0.15%	0.011%
64	17.542	2422	2426	2430	rBV3	5790	10143	1.08%	0.084%
65	17.601	2432	2436	2442	rBV2	3536	4679	0.50%	0.039%
66	17.689	2444	2451	2458	rVV	515592	784721	83.56%	6.473%
67	17.742	2458	2460	2461	rVV2	3336	2504	0.27%	0.021%
68	17.789	2461	2468	2474	rVV2	486611	720381	76.70%	5.942%
69	17.853	2477	2479	2485	rVB3	4135	4697	0.50%	0.039%
70	17.941	2487	2494	2497	rBV3	2839	5314	0.57%	0.044%
71	18.035	2505	2510	2511	rBV	1359	1974	0.21%	0.016%
72	18.171	2530	2533	2536	rBV2	1254	1696	0.18%	0.014%
73	18.270	2546	2550	2551	rBV2	1262	1665	0.18%	0.014%
74	18.329	2555	2560	2561	rBV2	5322	6053	0.64%	0.050%
75	18.376	2564	2568	2569	rVB2	1572	1541	0.16%	0.013%
76	18.417	2569	2575	2580	rBV6	7811	14960	1.59%	0.123%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

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 >
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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Title : SVOA CALIBRATION

77	18.482	2582	2586	2590	rBV4	26146	34426	3.67%	0.284%
78	18.535	2590	2595	2603	rBV	106171	139721	14.88%	1.152%
79	18.629	2609	2611	2615	rVB4	5776	5559	0.59%	0.046%
80	18.740	2626	2630	2636	rVB3	14971	18096	1.93%	0.149%
81	18.805	2636	2641	2648	rVV	23418	38153	4.06%	0.315%
82	18.870	2650	2652	2654	rVB3	3823	2360	0.25%	0.019%
83	18.964	2663	2668	2675	rBV8	14246	26202	2.79%	0.216%
84	19.058	2679	2684	2689	rVB3	26223	39567	4.21%	0.326%
85	19.128	2691	2696	2702	rVB2	64498	83756	8.92%	0.691%
86	19.181	2702	2705	2706	rBV2	6541	5784	0.62%	0.048%
87	19.252	2716	2717	2718	rVB	4792	1689	0.18%	0.014%
88	19.393	2738	2741	2745	rBV4	4045	5131	0.55%	0.042%
89	19.469	2749	2754	2759	rBV	46623	67152	7.15%	0.554%
90	19.675	2786	2789	2794	rBV4	16907	29600	3.15%	0.244%
91	19.792	2805	2809	2813	rBV3	46883	67079	7.14%	0.553%
92	20.051	2848	2853	2862	rVB	666688	939165	100.00%	7.747%
93	20.186	2871	2876	2882	rBV	659790	803944	85.60%	6.631%
94	20.309	2894	2897	2901	rBV4	18516	26091	2.78%	0.215%
95	20.538	2931	2936	2942	rBV	84442	116743	12.43%	0.963%
96	20.909	2994	2999	3008	rBV5	56497	117214	12.48%	0.967%
97	21.578	3109	3113	3126	rVB4	61333	118812	12.65%	0.980%
98	21.990	3176	3183	3189	rVB	503197	877087	93.39%	7.235%
99	25.221	3723	3733	3744	rVB	312644	889292	94.69%	7.335%
100	25.456	3765	3773	3786	rVB2	310369	915131	97.44%	7.548%

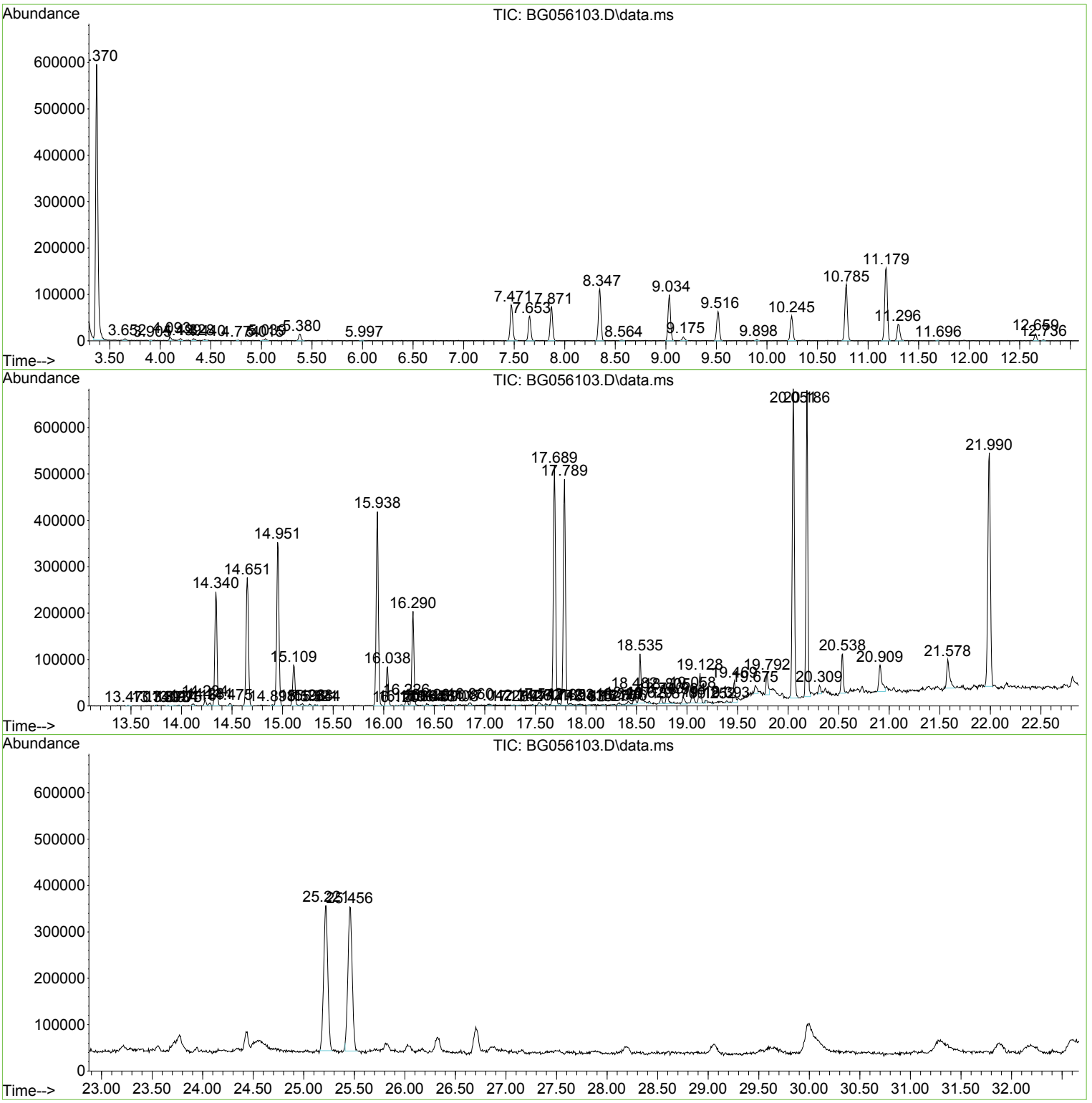
Sum of corrected areas: 12123517

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

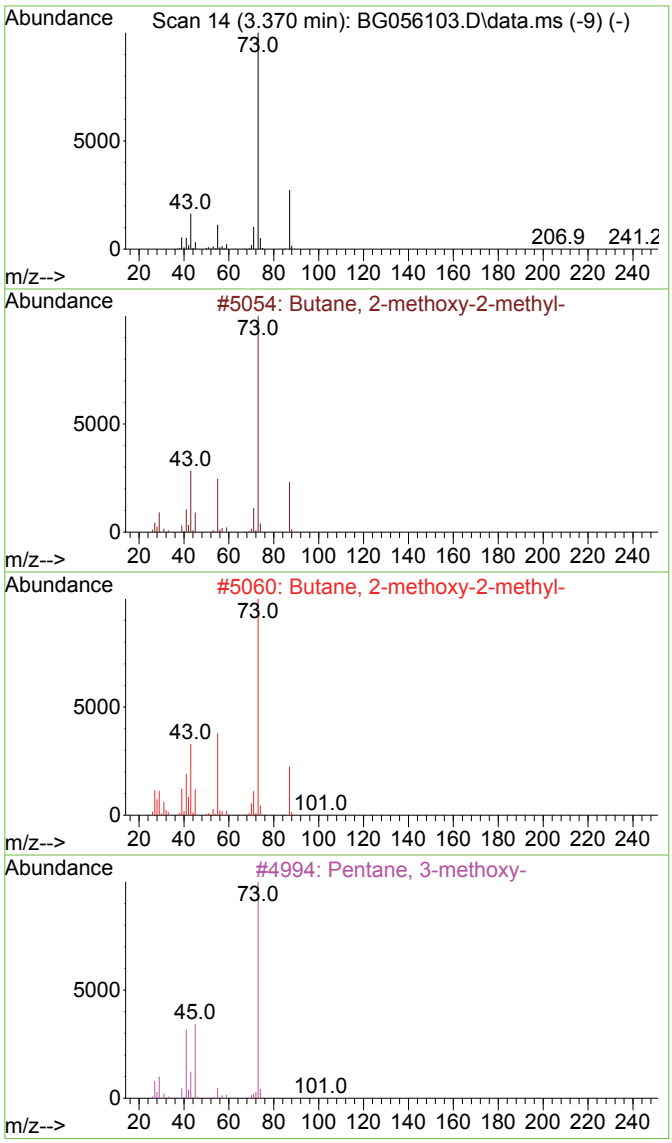
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

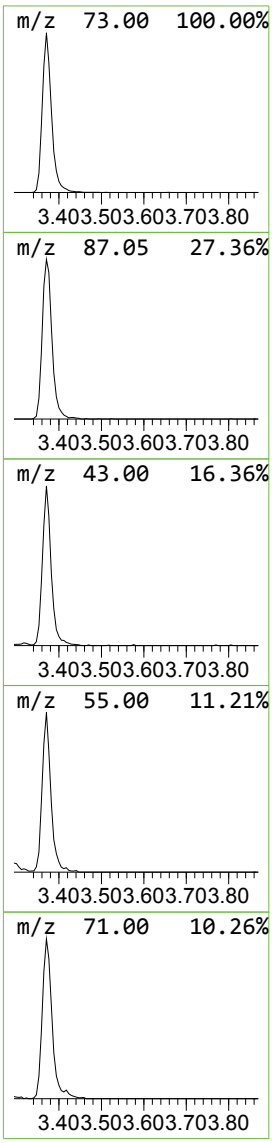
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.370	91.53 ng/ul	889907	1,4-Dichlorobenzene-d4	8.341

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	9
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
4	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	7





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

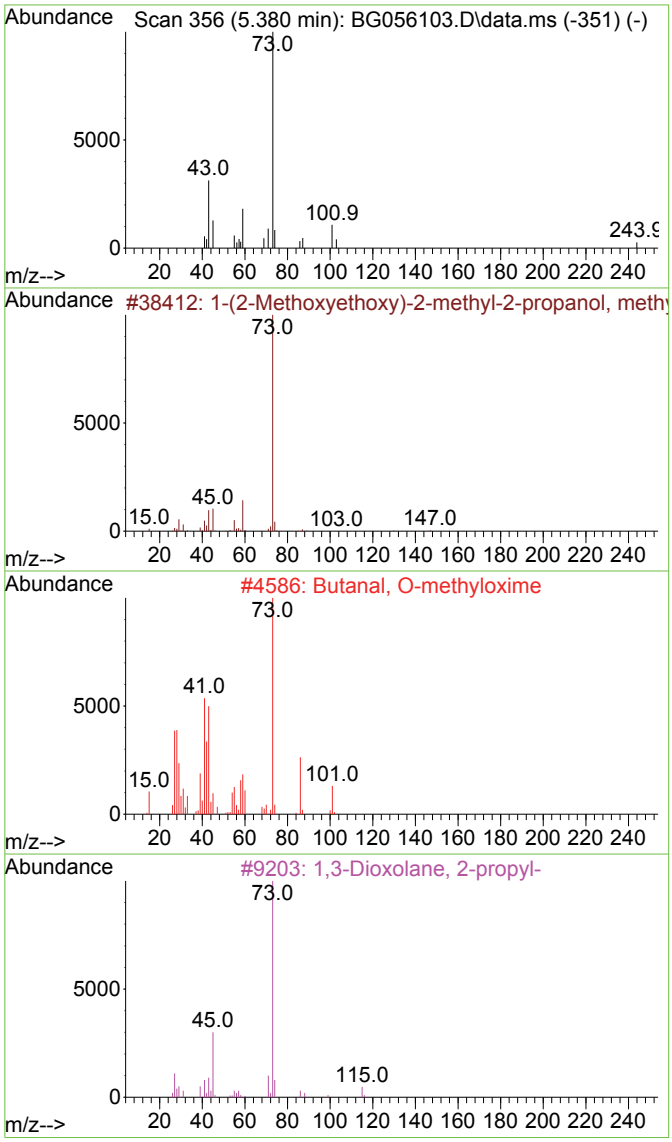
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

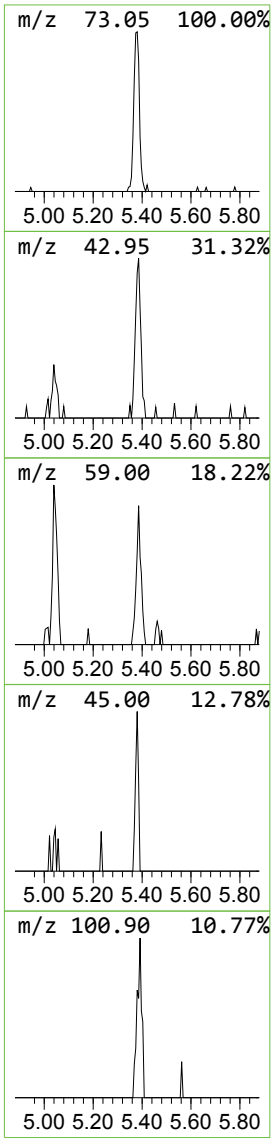
****

Peak Number 2 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.380	2.46 ng/ul	23896	1,4-Dichlorobenzene-d4	8.341

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	38
2	Butanal, O-methyloxime	101	C5H11NO	031376-98-4	33
3	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	25
4	Silane, trimethyl(3-methylbutyl)-	144	C8H20Si	018291-15-1	9
5	3-Heptanol, 2,4-dimethyl-	144	C9H20O	019549-72-5	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

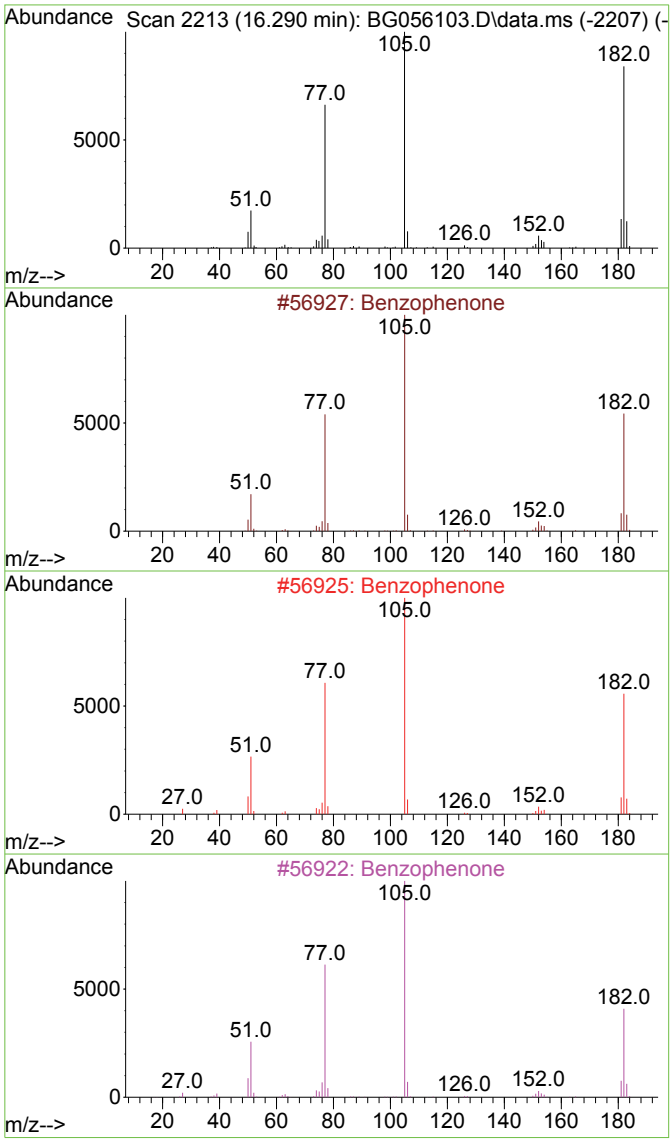
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

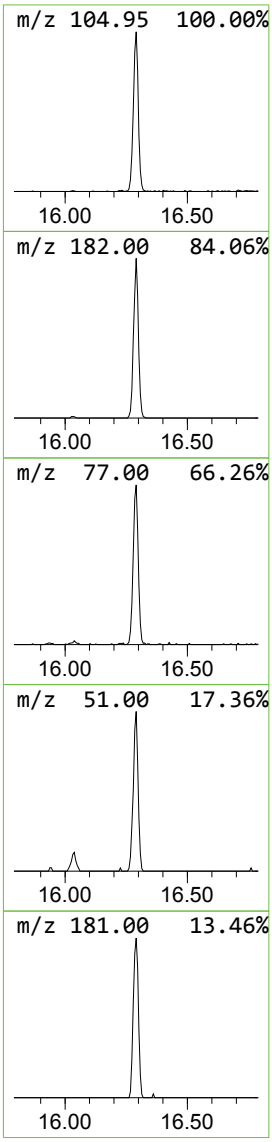
****

Peak Number 3 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.290	10.38 ng/ul	287269	Acenaphthene-d10	14.951

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	94
3	Benzophenone	182	C13H10O	000119-61-9	94
4	Benzophenone	182	C13H10O	000119-61-9	90
5	Benzenamine, N-(3-pyridinylmethy...	182	C12H10N2	029722-97-2	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

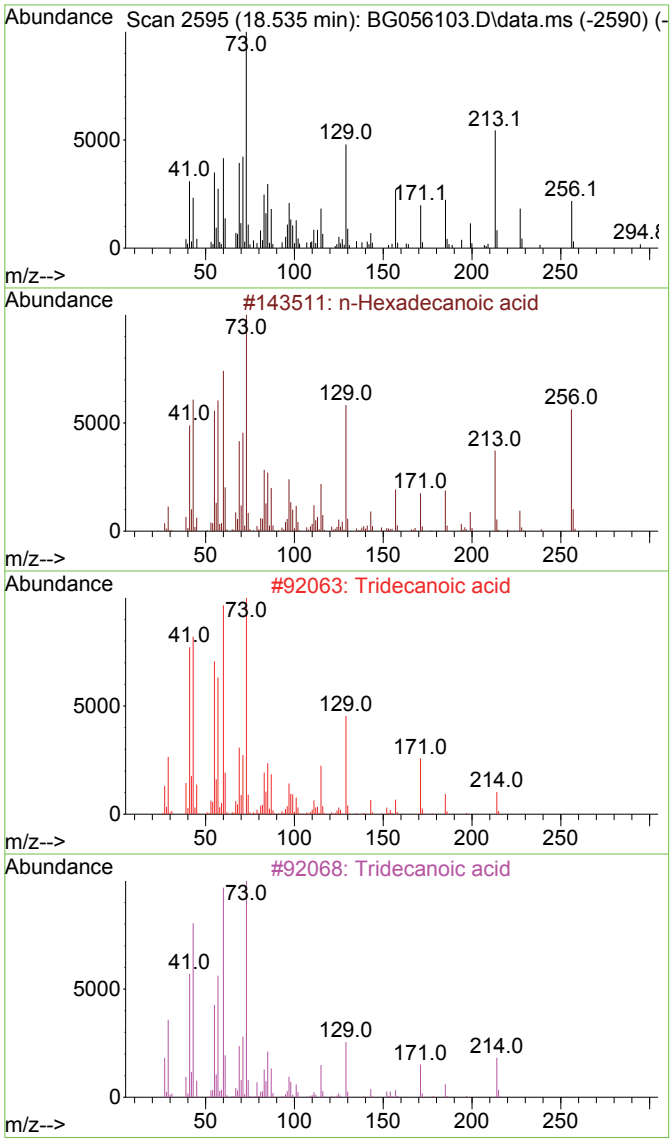
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

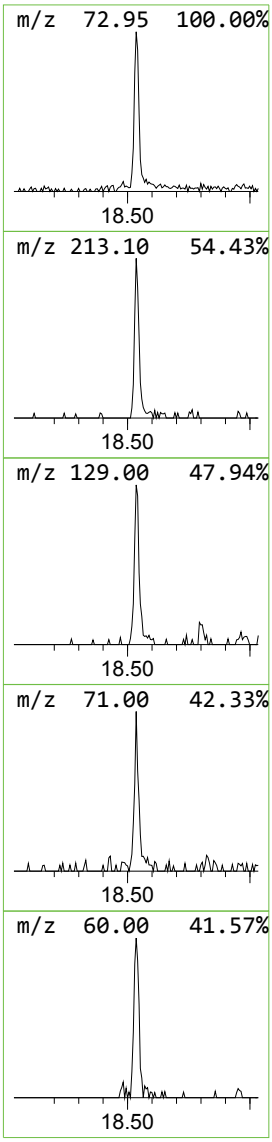
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.535	3.56 ng/ul	139721	Phenanthrene-d10	17.689

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70
2	Tridecanoic acid	214	C13H26O2	000638-53-9	56
3	Tridecanoic acid	214	C13H26O2	000638-53-9	50
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	38
5	Tridecanoic acid	214	C13H26O2	000638-53-9	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

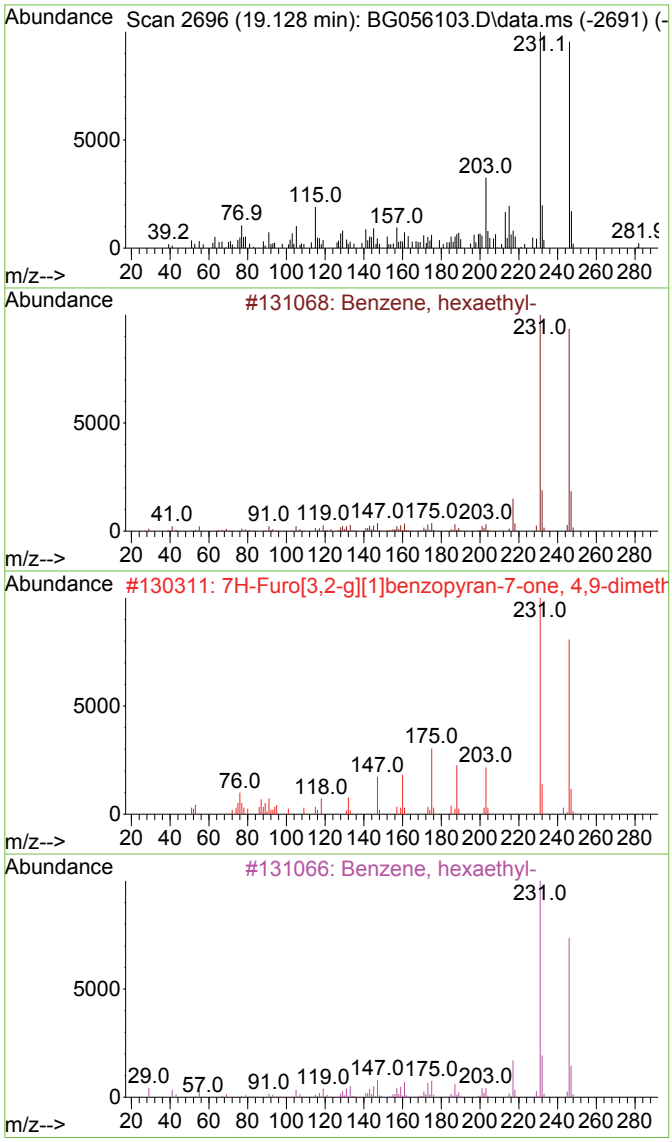
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

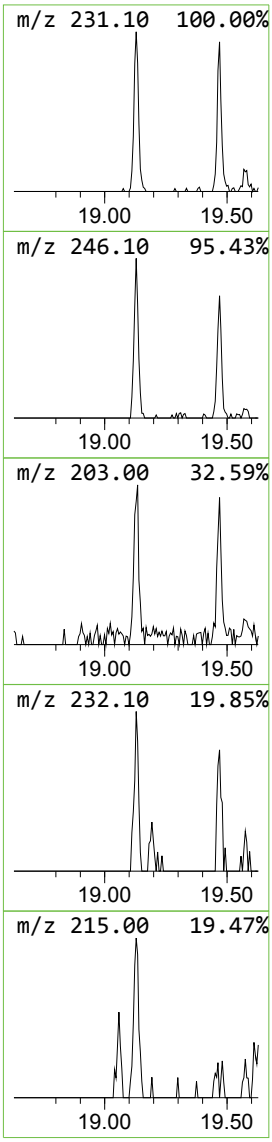
****

Peak Number 5 Benzene, hexaethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.128	2.13 ng/ul	83756	Phenanthrene-d10	17.689

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, hexaethyl-	246	C18H30	000604-88-6	76
2	7H-Furo[3,2-g][1]benzopyran-7-on...	246	C13H10O5	000482-27-9	72
3	Benzene, hexaethyl-	246	C18H30	000604-88-6	68
4	Pimpinellin	246	C13H10O5	000131-12-4	64
5	Benzene, 1,1'-thiobis[4-methoxy-	246	C14H14O2S	003393-77-9	49





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Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

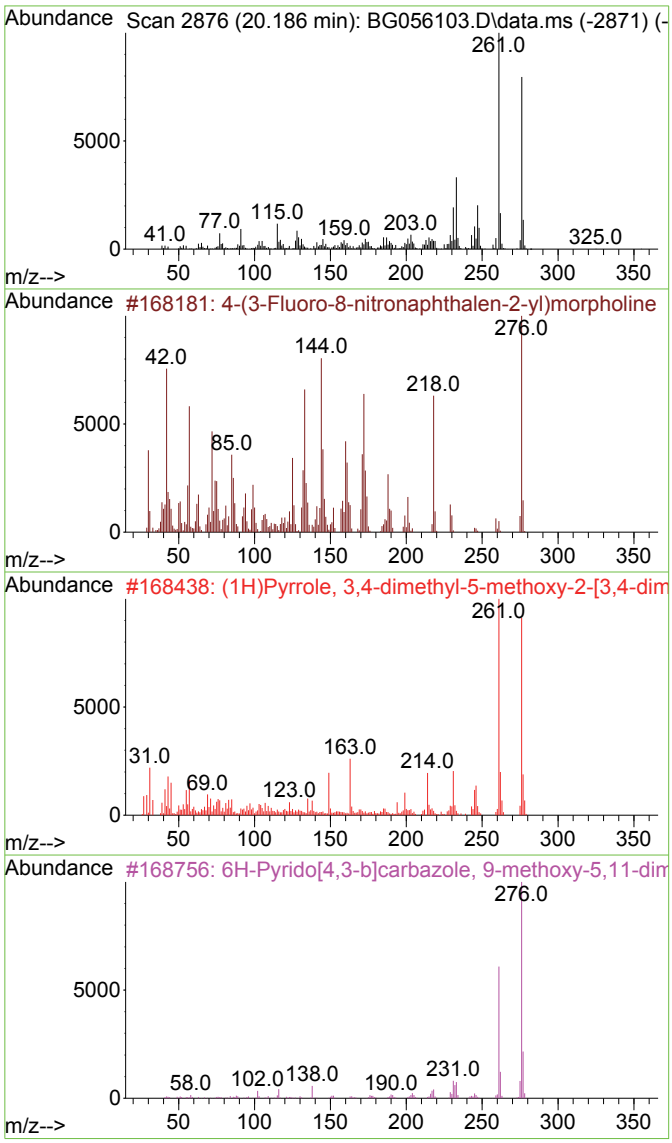
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

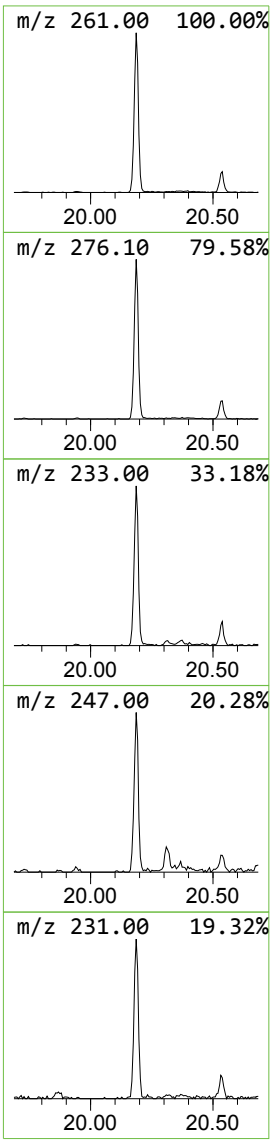
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4-(3-Fluoro-8-nitronaphthal... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.186	18.33 ng/ul	803944	Chrysene-d12	21.990

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(3-Fluoro-8-nitronaphthalen-2-...	276	C14H13FN2O3	1000409-37-1	91
2	(1H)Pyrrole, 3,4-dimethyl-5-meth...	276	C15H20N2OS	1000197-22-6	74
3	6H-Pyrido[4,3-b]carbazole, 9-met...	276	C18H16N2O	010371-86-5	72
4	3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	70
5	5H-Furo[3,2-g][1]benzopyran-5-on...	276	C14H12O6	000668-10-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
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Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

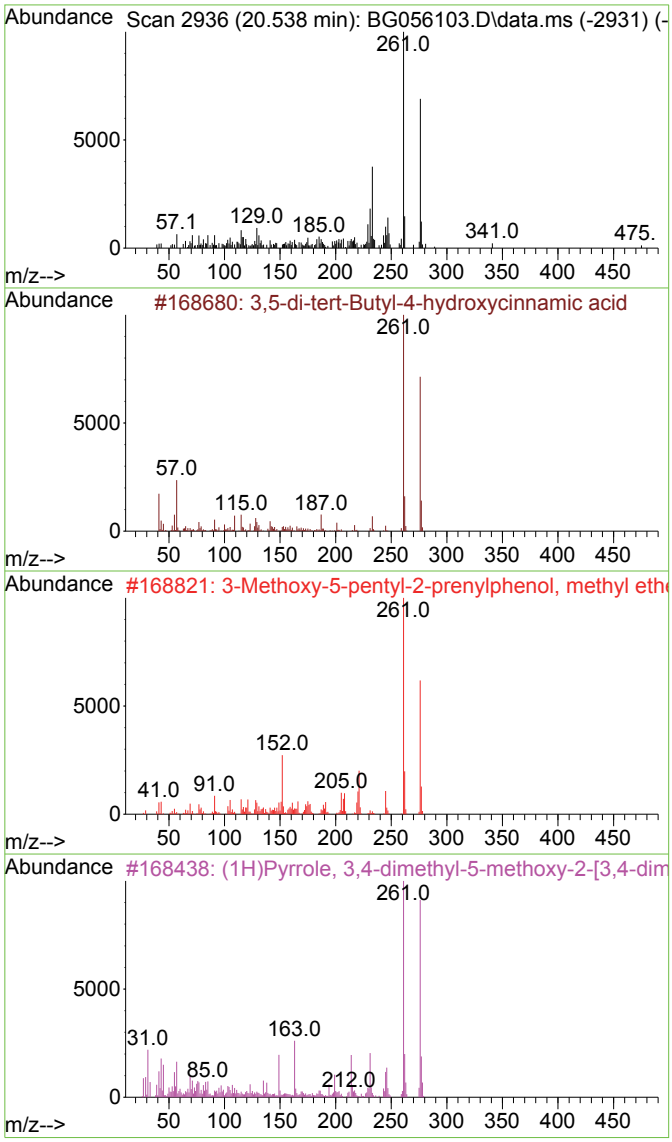
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

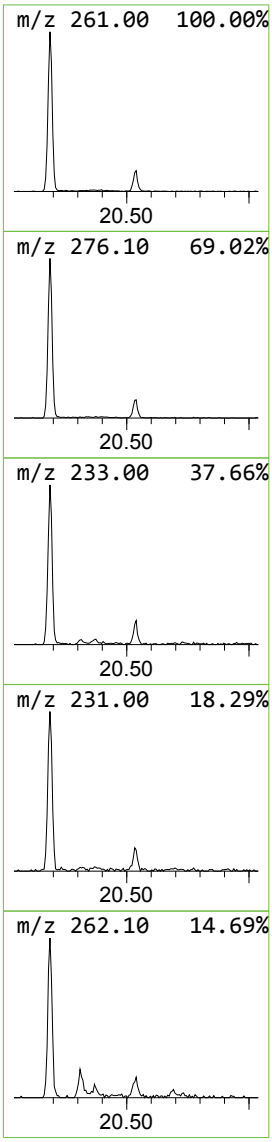
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.538	2.66 ng/ul	116743	Chrysene-d12	21.990

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	68
2	3-Methoxy-5-pentyl-2-prenylpheno...	276	C18H28O2	083036-53-7	64
3	(1H)Pyrrole, 3,4-dimethyl-5-meth...	276	C15H20N2OS	1000197-22-6	59
4	5H-Furo[3,2-g][1]benzopyran-5-on...	276	C14H12O6	000668-10-0	53
5	4-Ethyl-7-iodotropolone	276	C9H9IO2	004508-96-7	47





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Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

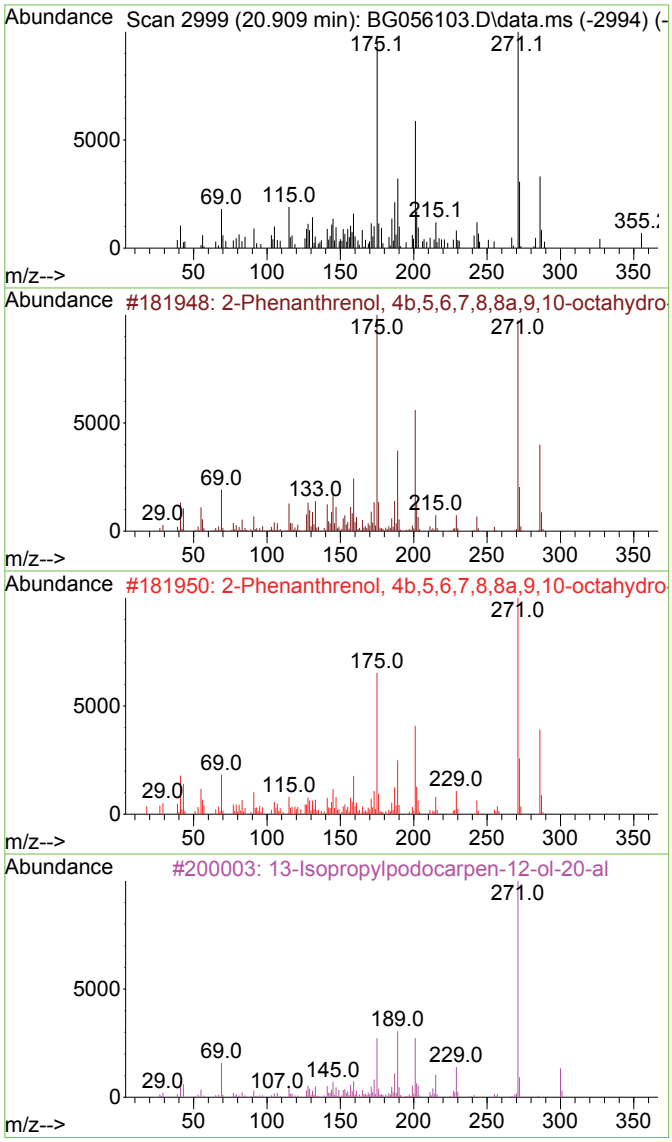
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

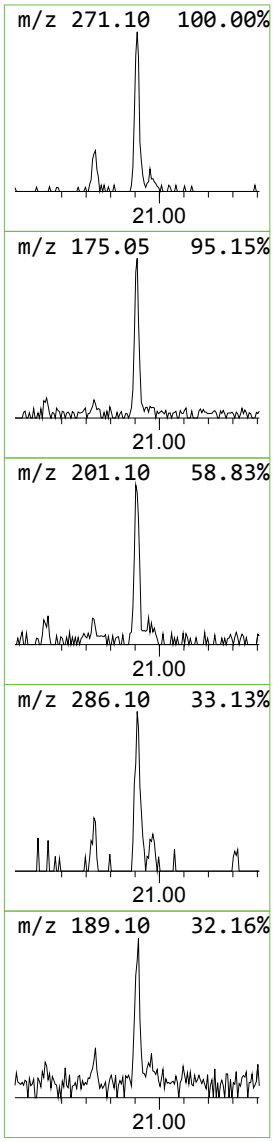
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.909	2.67 ng/ul	117214	Chrysene-d12	21.990

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	96
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	92
3	13-Isopropylpodocarpen-12-ol-20-al	300	C20H28O2	024035-37-8	43
4	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	38
5	Crinan-1-ol, 2,3-didehydro-,(1.a...	271	C16H17NO3	1000111-31-3	30





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Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

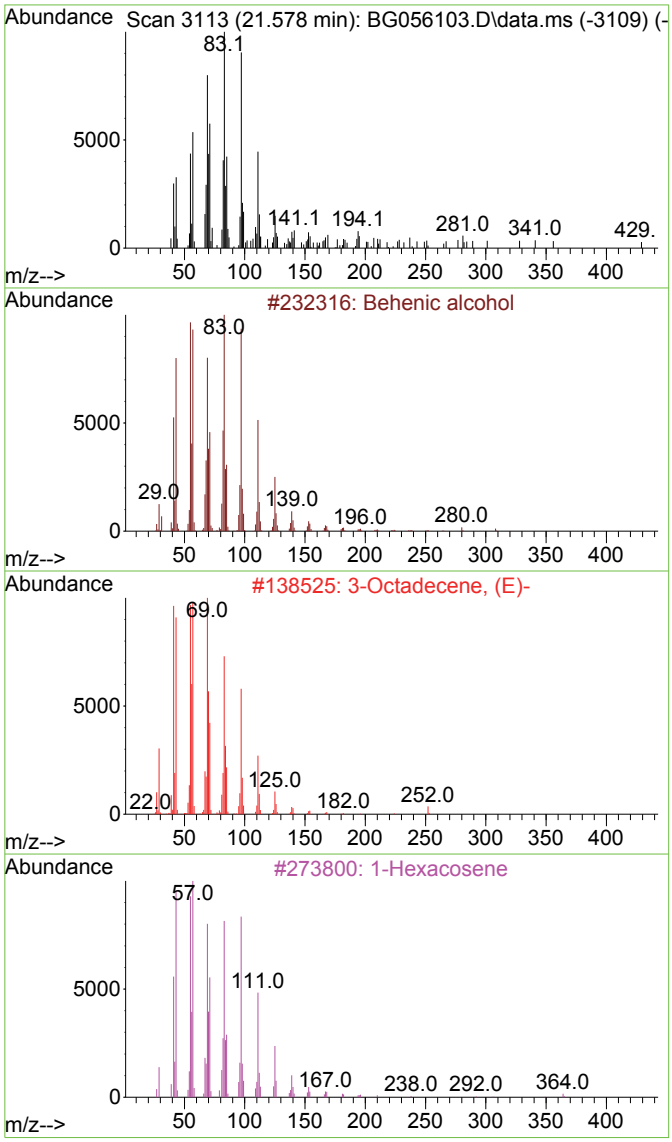
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

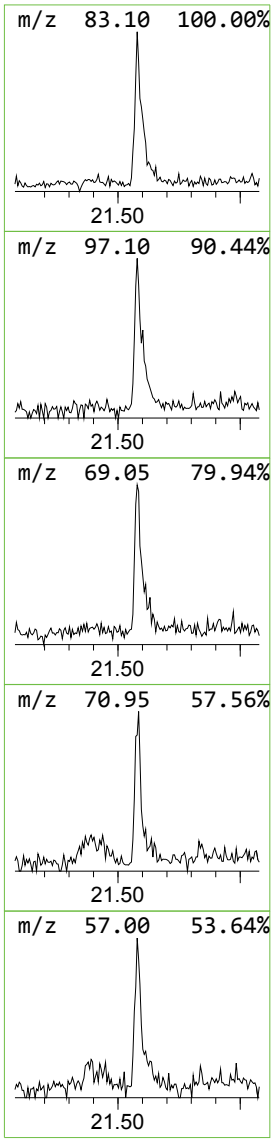
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Behenic alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.578	2.71 ng/ul	118812	Chrysene-d12	21.990

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	90
2	3-Octadecene, (E)-	252	C18H36	007206-19-1	89
3	1-Hexacosene	364	C26H52	018835-33-1	87
4	1-Hexadecanol	242	C16H34O	036653-82-4	87
5	Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYF1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	3.370	91.5	ng/ul	889907	1	8.341	194443	20.0
unknown-01	5.380	2.5	ng/ul	23896	1	8.341	194443	20.0
Benzophenone	16.290	10.4	ng/ul	287269	3	14.951	553608	20.0
n-Hexadecanoic ...	18.535	3.6	ng/ul	139721	4	17.689	784721	20.0
Benzene, hexaet...	19.128	2.1	ng/ul	83756	4	17.689	784721	20.0
4-(3-Fluoro-8-n...	20.186	18.3	ng/ul	803944	5	21.990	877087	20.0
3,5-di-tert-But...	20.538	2.7	ng/ul	116743	5	21.990	877087	20.0
2-Phenanthrenol...	20.909	2.7	ng/ul	117214	5	21.990	877087	20.0
Behenic alcohol	21.578	2.7	ng/ul	118812	5	21.990	877087	20.0