LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056123.D
Acq On : 25 Dec 2022 16:12
Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

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: BG056123.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.372	9	14	28	rVB	635673	963311	86.86%	6.709%
2	3.648	57	61	66	rVV2	5207	6776	0.61%	0.047%
3	4.200	149	155	159	rVB	6832	9582	0.86%	0.067%
4	4.329	172	177	183	rVB2	5265	8246	0.74%	0.057%
5	4.435	190	195	201	rVB4	2458	4697	0.42%	0.033%
6	5.381	350	356	363	rVB3	16895	28295	2.55%	0.197%
7	7.015	631	634	639	rVB2	4325	5221	0.47%	0.036%
8	7.473	699	712	719	rBV	95554	173437	15.64%	1.208%
9	7.655	735	743	749	rBV	133094	227407	20.51%	1.584%
10	7.867	773	779	786	rVV	100837	177685	16.02%	1.237%
11	8.342	853	860	870	rVB	118727	209792	18.92%	1.461%
12	9.036	970	978	988	rBV	155732	267757	24.14%	1.865%
13	9.171	995	1001	1009	rBV	11451	22225	2.00%	0.155%
14	9.518	1051	1060	1068	rBV	179533	308991	27.86%	2.152%
15	9.894	1120	1124	1129	rVB	4637	6727	0.61%	0.047%
16	10.240	1175	1183	1192	rBV	157094	286700	25.85%	1.997%
17	10.781	1268	1275	1283	rVB	165536	291829	26.31%	2.032%
18	11.180	1334	1343	1353	rBV	176707	328429	29.61%	2.287%
19	11.298	1356	1363	1371	rBV	101479	185116	16.69%	1.289%
20	13.818	1789	1792	1795	rBV2	2910	4140	0.37%	0.029%
21	13.865	1798	1800	1808	rVB2	3230	5236	0.47%	0.036%
22	14.283	1865	1871	1875	rVB2	17793	25811	2.33%	0.180%
23	14.341	1875	1881	1887	rBV	581460	838180	75.58%	5.837%
24	14.653	1926	1934	1941	rBV	352206	557619	50.28%	3.884%
25	14.788	1953	1957	1959	rBV2	3810	4718	0.43%	0.033%
26	14.905	1974	1977	1979	rBV2	5438	7119	0.64%	0.050%
27	14.952	1979	1985	1991	rVV	357789	557248	50.25%	3.881%
28	15.011	1991	1995	1998	rVV4	10667	15276	1.38%	0.106%
29	15.117	2006	2013	2020	rBV	115045	187359	16.89%	1.305%
30	15.187	2020	2025	2034	rVB3	20752	33577	3.03%	0.234%
31	15.264	2036	2038	2043	rVB2	3461	4160	0.38%	0.029%
32	15.328	2043	2049	2052	rBV3	2942	5473	0.49%	0.038%
33	15.845	2134	2137	2143	rVB3	14946	20664	1.86%	0.144%
34	15.939	2146	2153	2160	rBV	708291	1109003	100.00%	7.724%
35	15.998	2160	2163	2164	rVV2	3374	4241	0.38%	0.030%
36	16.039	2164	2170	2180	rVB	211337	316549	28.54%	2.205%

LSC Area Percent Report

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Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

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	16.174	2188	2193	2199	rVB4	7193	10648	0.96%	0.074%
38	16.292	2207	2213	2220	rVB	36961	54647	4.93%	0.381%
39	16.374	2222	2227	2228	rBV4	6303	9365	0.84%	0.065%
40	16.392	2228	2230	2238	rVB7	7524	10211	0.92%	0.071%
41	16.498	2243	2248	2252	rBV4	14627	19463	1.75%	0.136%
42	16.633	2263	2271	2276	rBV7	3399	7203	0.65%	0.050%
43	16.715	2282	2285	2288	rVB2	4626	4871	0.44%	0.034%
44	17.044	2338	2341	2344	rBV3	4630	6410	0.58%	0.045%
45	17.179	2361	2364	2368	rVB4	5565	7742	0.70%	0.054%
46	17.220	2368	2371	2374	rBV3	3325	4283	0.39%	0.030%
47	17.338	2386	2391	2393	rBV3	4635	7374	0.66%	0.051%
48	17.408	2398	2403	2409	rVB4	4472	6838	0.62%	0.048%
49	17.485	2409	2416	2419	rBV4	3411	8115	0.73%	0.057%
50	17.543	2422	2426	2428	rBV4	6494	9008	0.81%	0.063%
51	17.608	2433	2437	2441	rVB3	5252	7024	0.63%	0.049%
52	17.690	2443	2451	2457	rBV	477190	720613	64.98%	5.019%
53	17.731	2457	2458	2461	rVV3	5191	4435	0.40%	0.031%
54	17.790	2461	2468	2475	rVB	485061	739988	66.73%	5.154%
55	18.325	2555	2559	2563	rBV5	9860	12015	1.08%	0.084%
56	18.366	2563	2566	2568	rVB4	4058	4132	0.37%	0.029%
57	18.425	2568	2576	2580	rBV5	8754	20359	1.84%	0.142%
58	18.478	2582	2585	2590	rBV4	10912	15226	1.37%	0.106%
59	18.536	2590	2595	2603	rBV2	115231	157933	14.24%	1.100%
60	18.613	2604	2608	2613	rVB3	7305	15470	1.39%	0.108%
61	18.742	2625	2630	2634	rVB7	7485	12814	1.16%	0.089%
62	18.807	2636	2641	2644	rBV3	9123	14019	1.26%	0.098%
63	18.901	2653	2657	2660	rBV5	6681	12414	1.12%	0.086%
64	18.954	2662	2666	2671	rVV6	14176	23123	2.09%	0.161%
65	19.048	2678	2682	2686	rVV3	15177	27738	2.50%	0.193%
66	19.089	2686	2689	2691	rVV3	13625	15642	1.41%	0.109%
67	19.130	2691	2696	2701	rVB5	29119	45080	4.06%	0.314%
68	19.335	2729	2731	2733	rBV3	7938	6831	0.62%	0.048%
69	19.471	2747	2754	2760	rBV8	25879	56678	5.11%	0.395%
70	19.570	2769	2771	2772	rBV2	8178	5976	0.54%	0.042%
71	19.647	2781	2784	2787	rBV4	9366	12008	1.08%	0.084%
72	19.788	2805	2808	2812	rBV3	42634	58908	5.31%	0.410%
73	19.946	2832	2835	2839	rBV6	11822	16087	1.45%	0.112%
74	20.052	2847	2853	2859	rBV	621249	893220	80.54%	6.221%
75	20.181	2871	2875	2880	rBV	172296	225606	20.34%	1.571%
76	20.317	2895	2898	2900	rBV4	8448	7344	0.66%	0.051%

LSC Area Percent Report

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Data File : BG056123.D
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Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Title : SVOA CALIBRATION

77	20.370	2905	2907	2912	rBV6	9776	11857	1.07%	0.083%
78	20.481	2924	2926	2928	rBV3	6774	4130	0.37%	0.029%
79	20.540	2933	2936	2938	rVV4	20953	25696	2.32%	0.179%
80	20.857	2988	2990	2991	rBV2	9817	6595	0.59%	0.046%
81	21.004	3012	3015	3018	rVB2	14802	21514	1.94%	0.150%
82	21.069	3024	3026	3030	rVB5	12681	12373	1.12%	0.086%
83	21.351	3070	3074	3077	rBV5	14544	18470	1.67%	0.129%
84	21.574	3108	3112	3125	rVB3	84562	157670	14.22%	1.098%
85	21.697	3131	3133	3138	rVB2	16925	16825	1.52%	0.117%
86	21.991	3176	3183	3193	rVB	492134	839682	75.72%	5.848%
87	22.814	3320	3323	3325	rBV3	18857	22147	2.00%	0.154%
88	23.554	3442	3449	3454	rBV10	22354	57252	5.16%	0.399%
89	24.429	3592	3598	3606	rVB3	44406	98103	8.85%	0.683%
90	25.217	3722	3732	3743	rVB	307746	879727	79.33%	6.127%
91	25.464	3764	3774	3784	rVB2	300813	865450	78.04%	6.027%
92	26.033	3868	3871	3878	rVB8	15114	29910	2.70%	0.208%
93	26.697	3976	3984	3994	rVB3	65164	213987	19.30%	1.490%
94	26.838	4000	4008	4026	rVB5	96852	359189	32.39%	2.502%
95	27.855	4174	4181	4182	rBV7	21894	41290	3.72%	0.288%
96	28.178	4228	4236	4237	rBV8	24616	52087	4.70%	0.363%
97	29.976	4534	4542	4543	rBV7	53450	102439	9.24%	0.713%
98	30.234	4584	4586	4587	rBV2	8805	5352	0.48%	0.037%
99	32.050	4893	4895	4897	rBV3	6821	5668	0.51%	0.039%
100	32.496	4969	4971	4973	rBV3	7993	5761	0.52%	0.040%

Sum of corrected areas: 14358631

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056123.D
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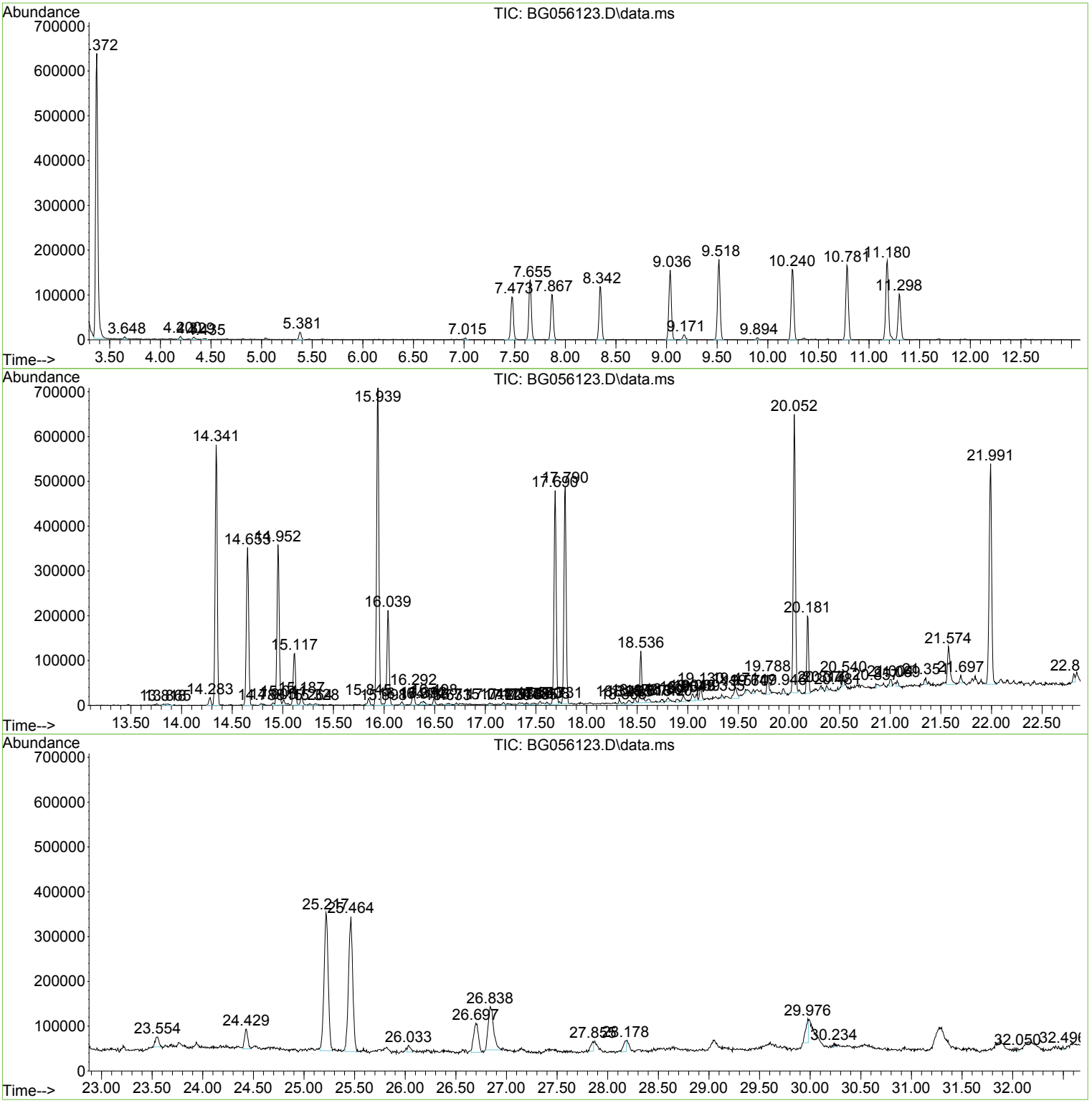
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

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 : BG056123.D
Acq On : 25 Dec 2022 16:12
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ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

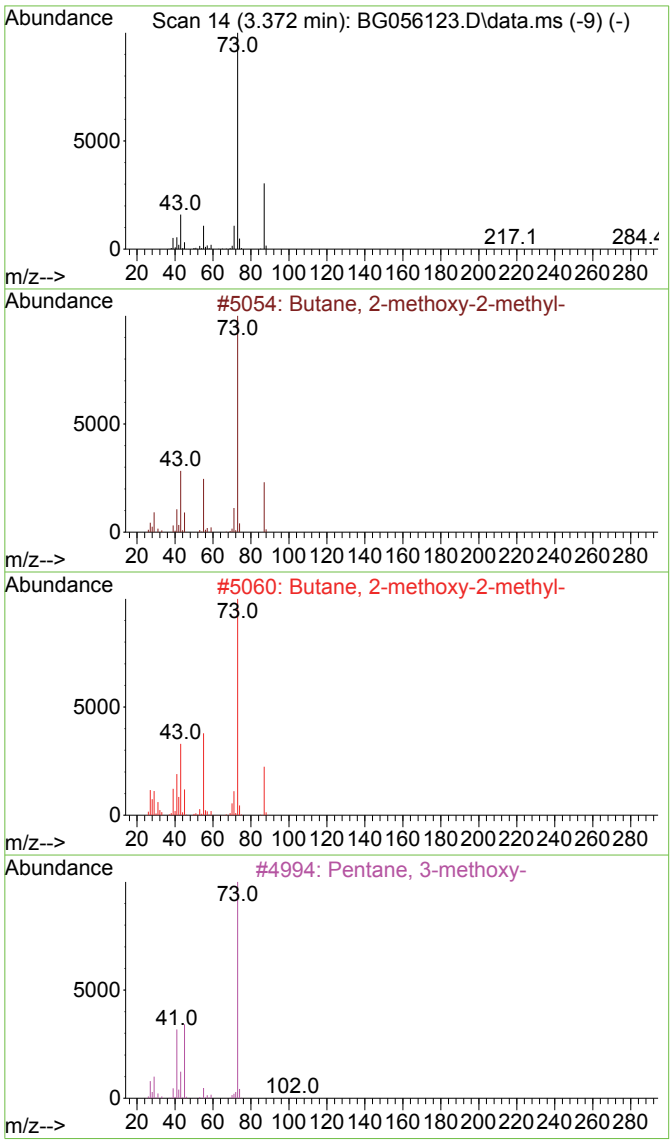
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

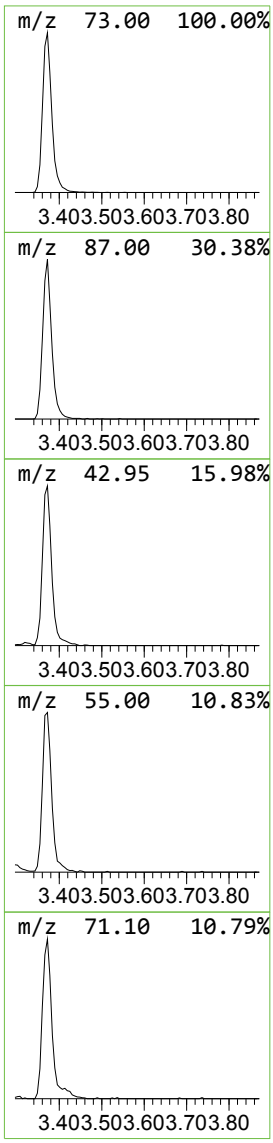
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.372	91.83 ng/ul	963311	1,4-Dichlorobenzene-d4	8.342

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	64
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
4	1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	9
5	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056123.D
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Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

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

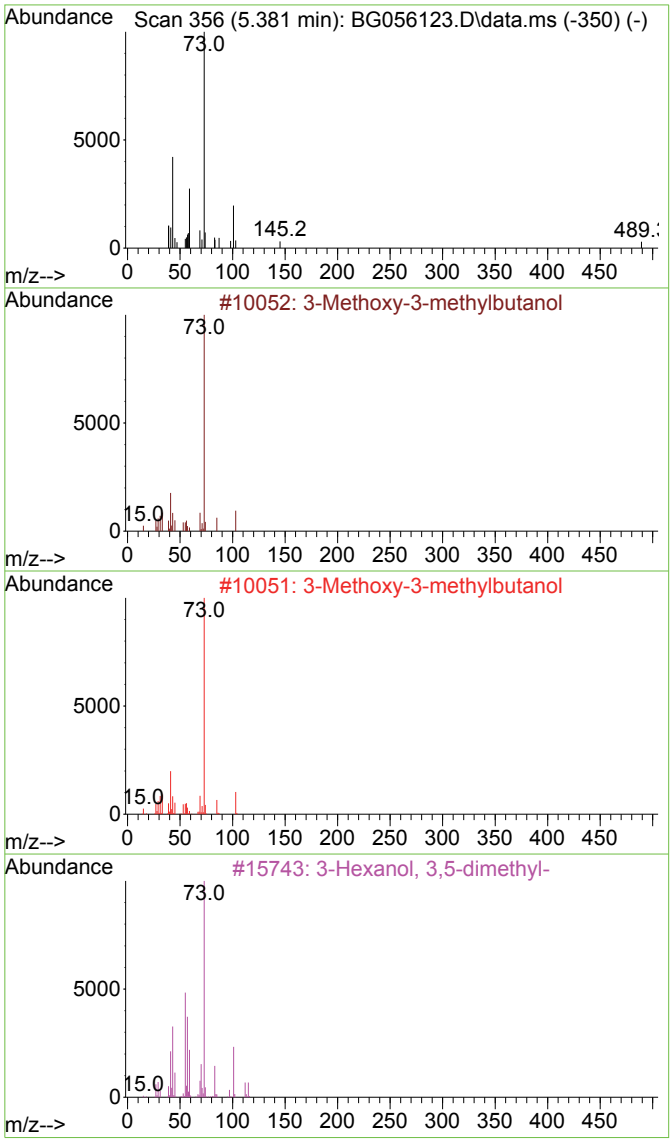
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

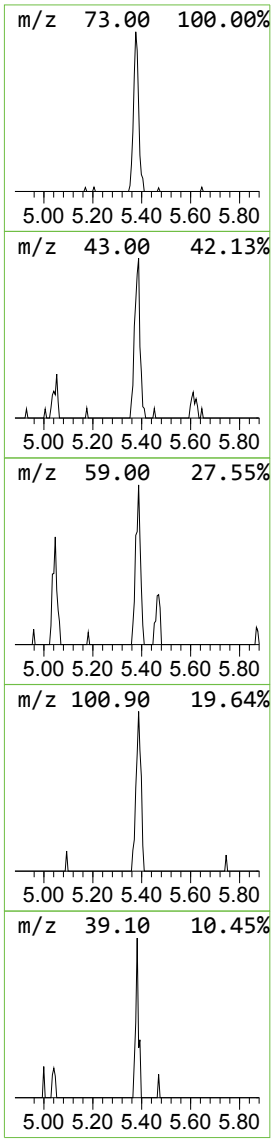
****

Peak Number 2 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.381	2.70 ng/ul	28295	1,4-Dichlorobenzene-d4	8.342

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methoxy-3-methylbutanol	118	C6H14O2	056539-66-3	10
2	3-Methoxy-3-methylbutanol	118	C6H14O2	056539-66-3	10
3	3-Hexanol, 3,5-dimethyl-	130	C8H18O	004209-91-0	9
4	1,3-Dioxolane, 2-butyl-	130	C7H14O2	004360-76-3	9
5	Isopropyl 5,11-dihydroxy-3,7,11-...	314	C18H34O4	1000223-34-1	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056123.D
Acq On : 25 Dec 2022 16:12
Operator : CG/JU
Sample : N6017-21
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ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

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

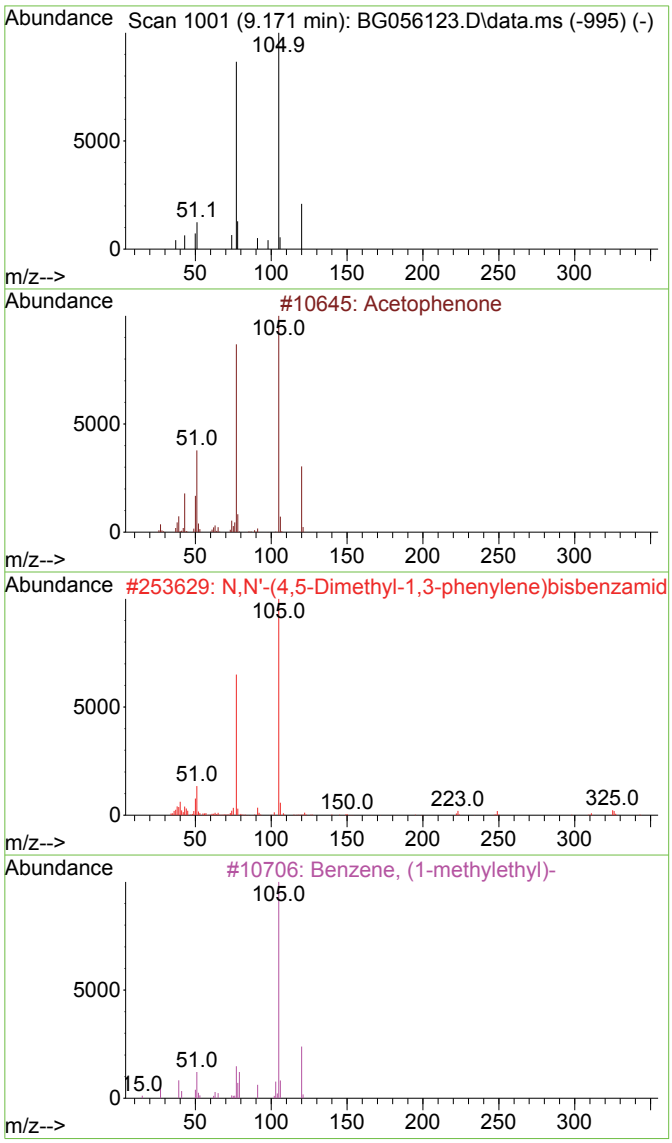
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

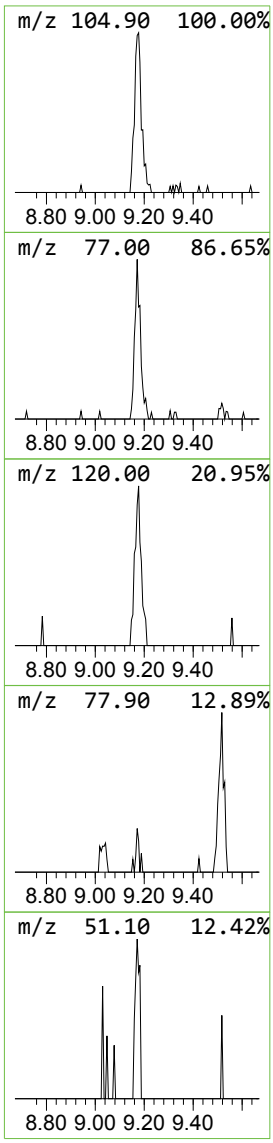
****

Peak Number 3 Aceto phenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.171	2.12 ng/ul	22225	1,4-Dichlorobenzene-d4	8.342

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone	120	C8H8O	000098-86-2	86
2	N,N'-(4,5-Dimethyl-1,3-phenylene...	344	C22H20N2O2	047507-95-9	45
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	25
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	16
5	1,5-Hexadiyne	78	C6H6	000628-16-0	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
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Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

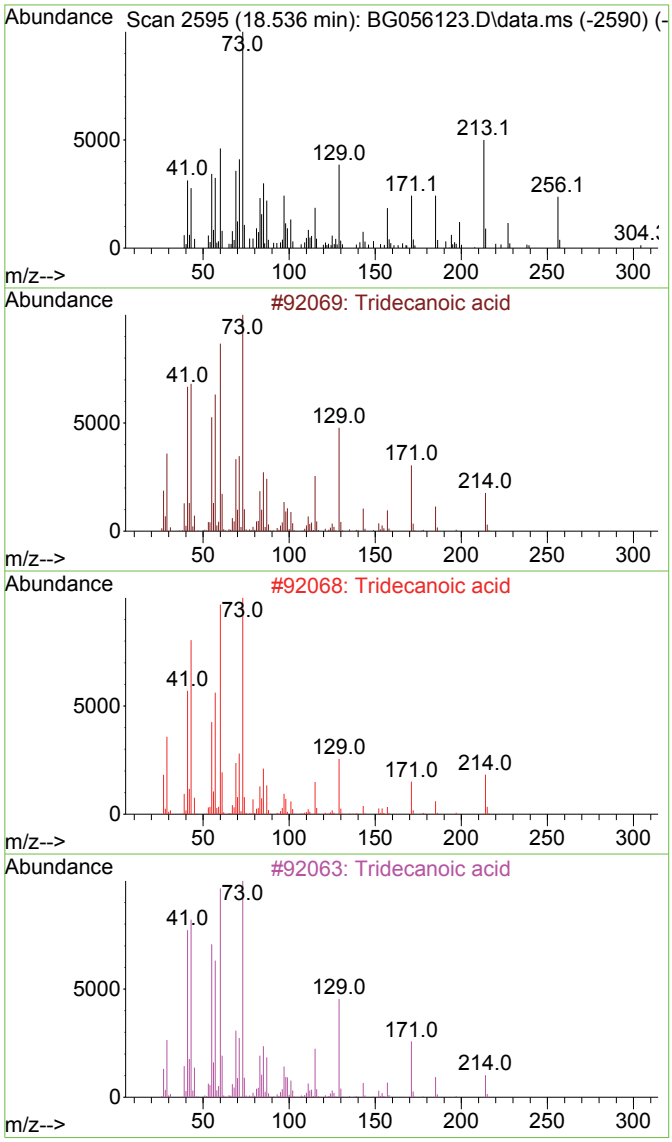
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

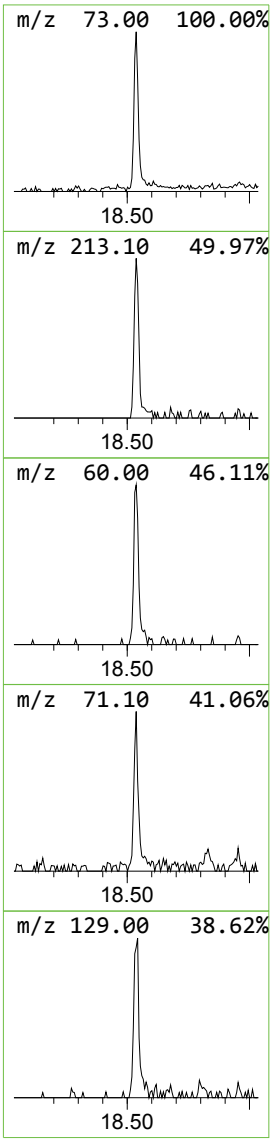
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.536	4.38 ng/ul	157933	Phenanthrene-d10	17.690

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	60
2	Tridecanoic acid	214	C13H26O2	000638-53-9	55
3	Tridecanoic acid	214	C13H26O2	000638-53-9	50
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	49
5	Tridecanoic acid	214	C13H26O2	000638-53-9	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056123.D
Acq On : 25 Dec 2022 16:12
Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

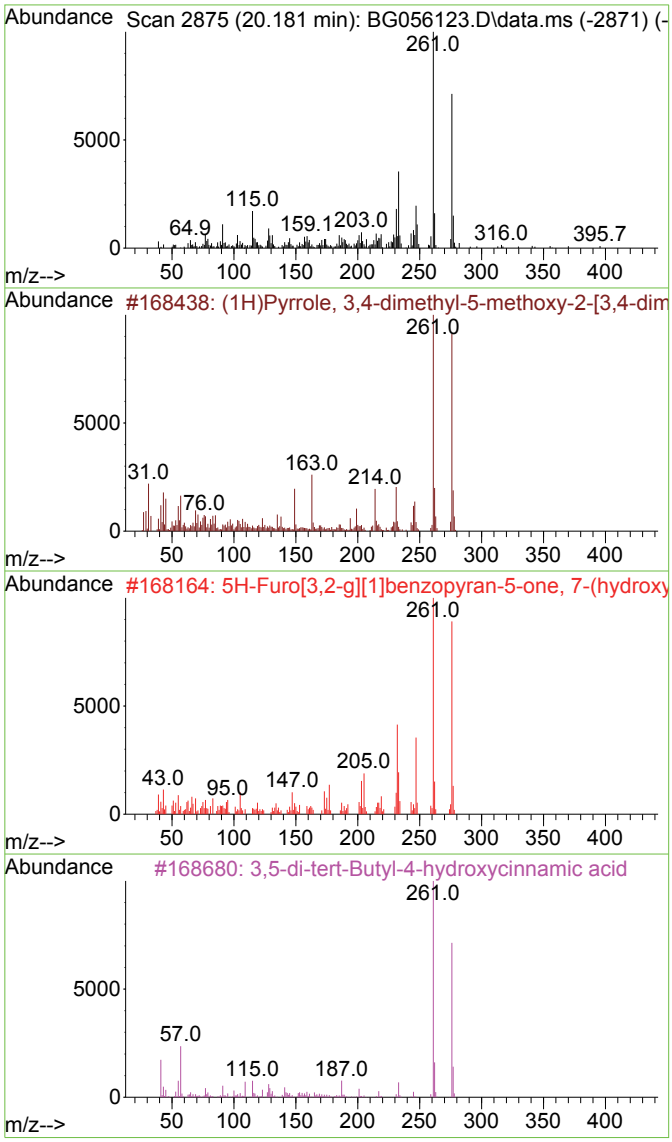
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

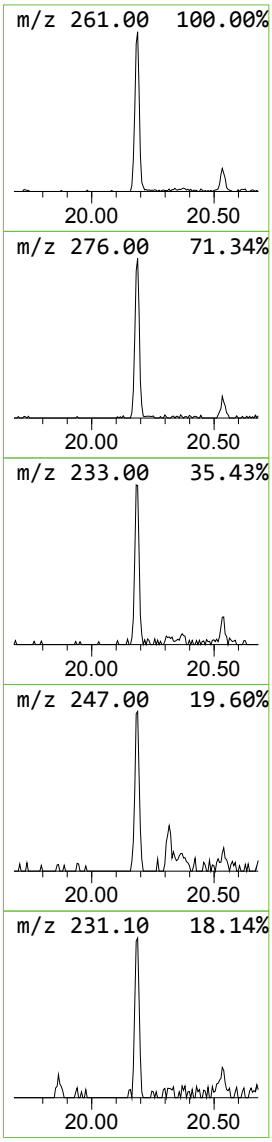
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (1H)Pyrrole, 3,4-dimethyl-5... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.182	5.37 ng/ul	225606	Chrysene-d12	21.991

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1H)Pyrrole, 3,4-dimethyl-5-meth...	276	C15H20N2OS	1000197-22-6	89
2	5H-Furo[3,2-g][1]benzopyran-5-on...	276	C14H12O6	000668-10-0	81
3	3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	70
4	5,6,7,8-Tetrahydrothieno[2,3-b]q...	276	C14H20N2SSi	1000496-40-4	64
5	4-(3-Fluoro-8-nitronaphthalen-2-...	276	C14H13FN2O3	1000409-37-1	56





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Data File : BG056123.D
Acq On : 25 Dec 2022 16:12
Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

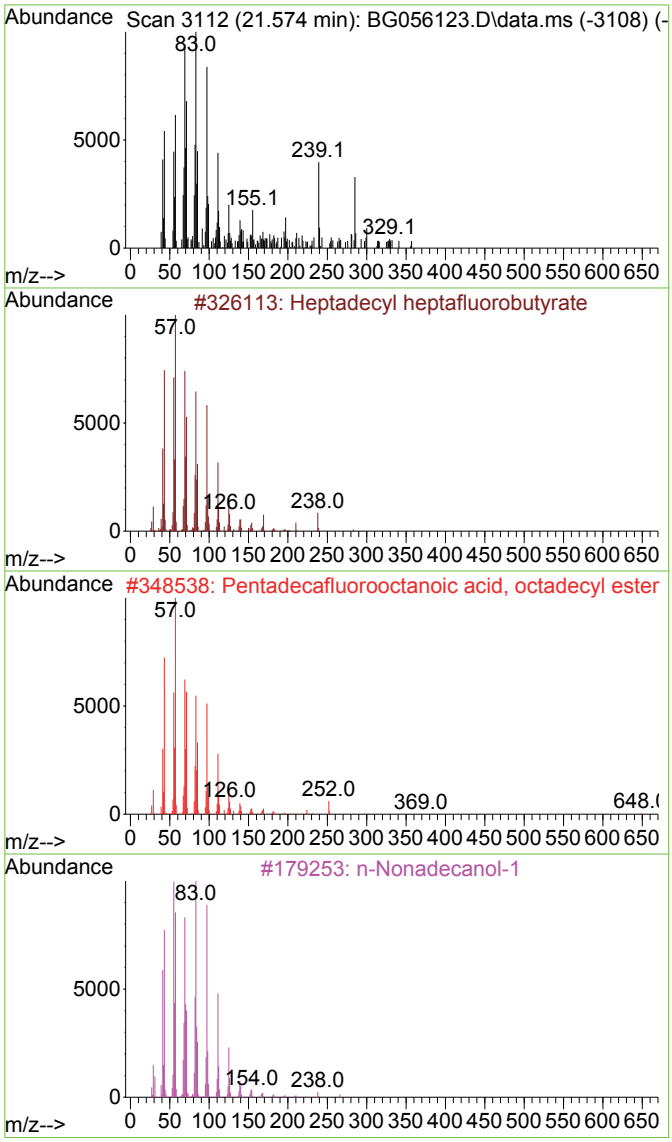
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

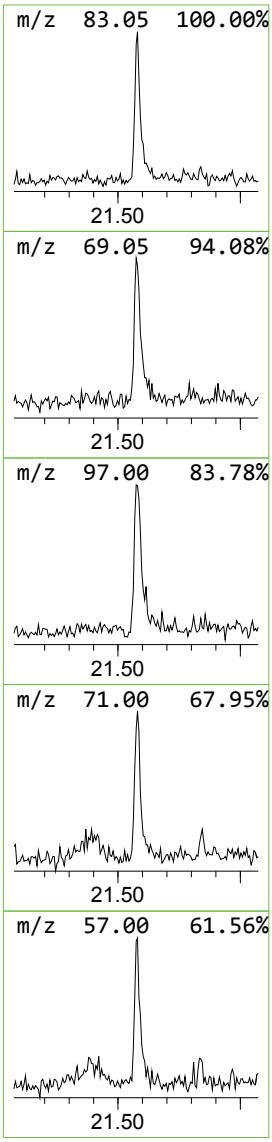
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Heptadecyl heptafluorobutyrate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.574	3.76 ng/ul	157670	Chrysene-d12	21.991

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	81
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	76
4	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	76
5	n-Pentadecanol	228	C15H32O	000629-76-5	76





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Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

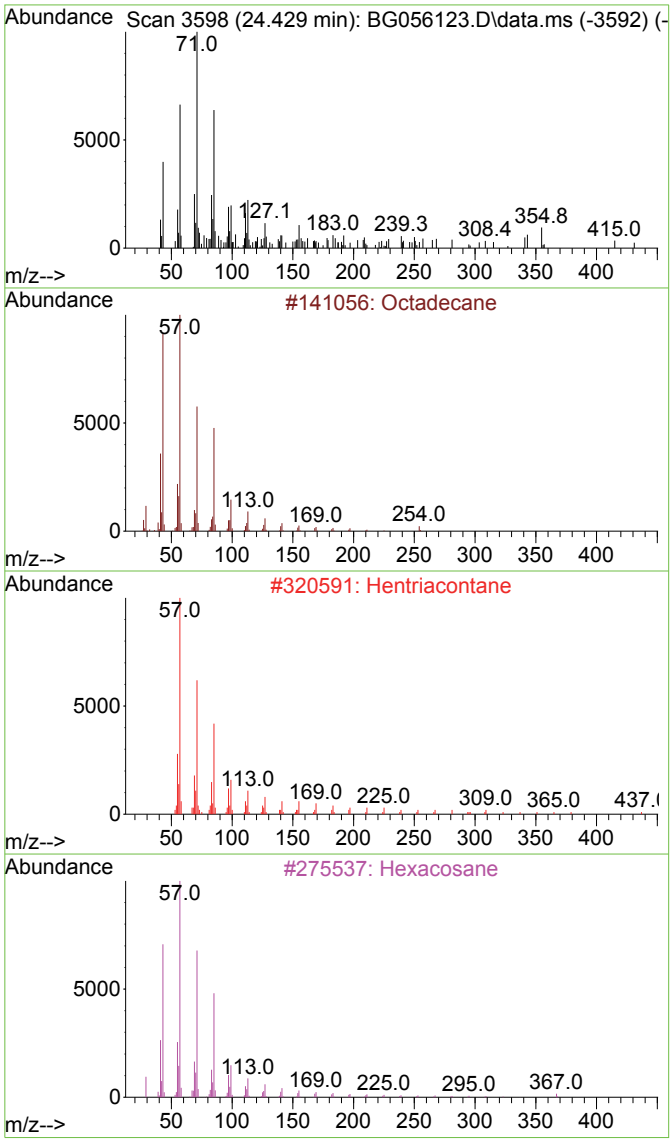
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

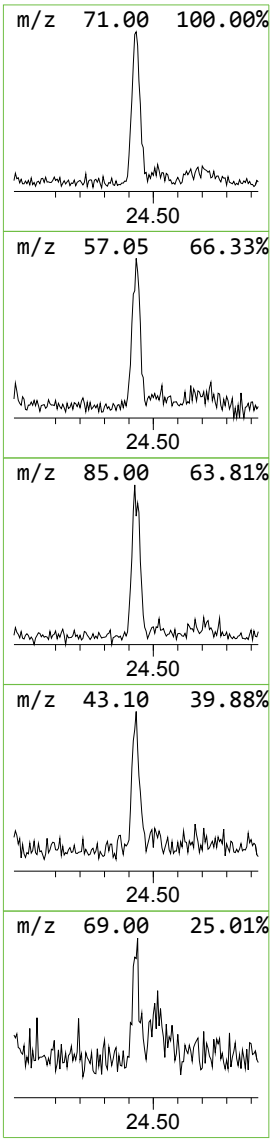
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.429	2.27 ng/ul	98103	Perylene-d12	25.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	76
2	Hentriacontane	437	C31H64	000630-04-6	74
3	Hexacosane	366	C26H54	000630-01-3	72
4	Heptacosane	380	C27H56	000593-49-7	72
5	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	72





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Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

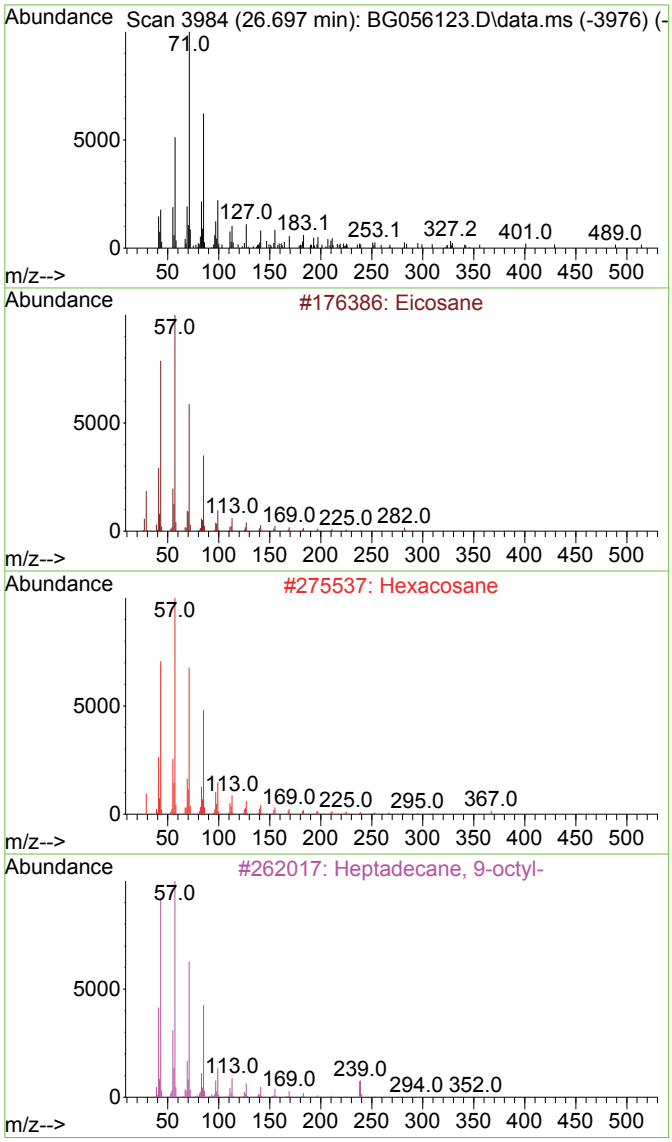
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

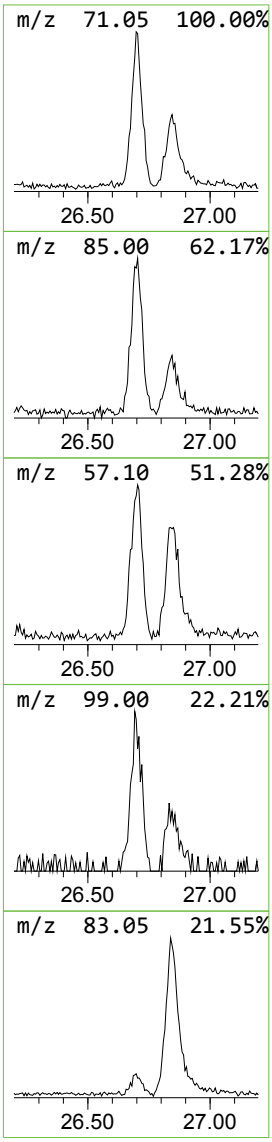
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.697	4.95 ng/ul	213987	Perylene-d12	25.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Hexacosane	366	C26H54	000630-01-3	87
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
4	Hexacosane	366	C26H54	000630-01-3	83
5	Hentriacontane	437	C31H64	000630-04-6	83





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Data File : BG056123.D
Acq On : 25 Dec 2022 16:12
Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

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

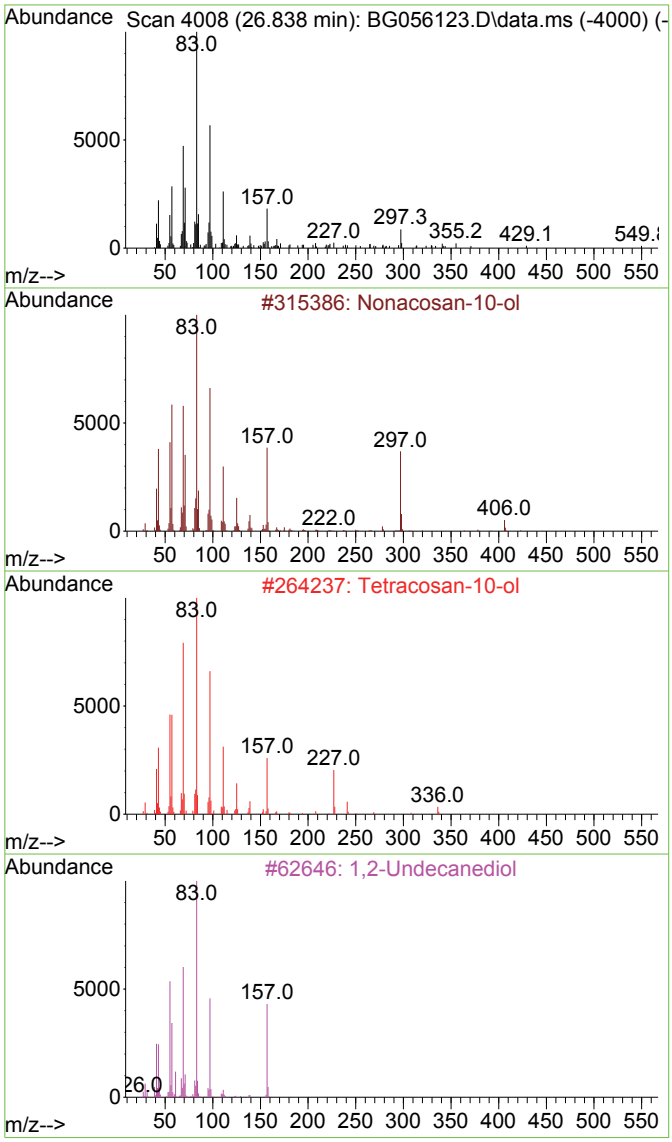
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

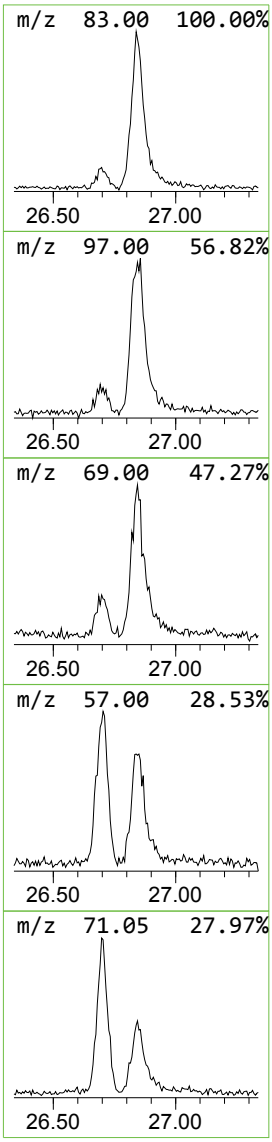
****

Peak Number 9 Nonacosan-10-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.838	8.30 ng/ul	359189	Perylene-d12	25.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonacosan-10-ol	424	C29H60O	000504-55-2	83
2	Tetracosan-10-ol	354	C24H50O	138966-95-7	64
3	1,2-Undecanediol	188	C11H24O2	013006-29-6	59
4	10-Nonadecanol	284	C19H40O	016840-84-9	59
5	1H,5H-Pyrrolo[1',2':3,4]imidazo[...	166	C10H18N2	054966-11-9	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056123.D
Acq On : 25 Dec 2022 16:12
Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

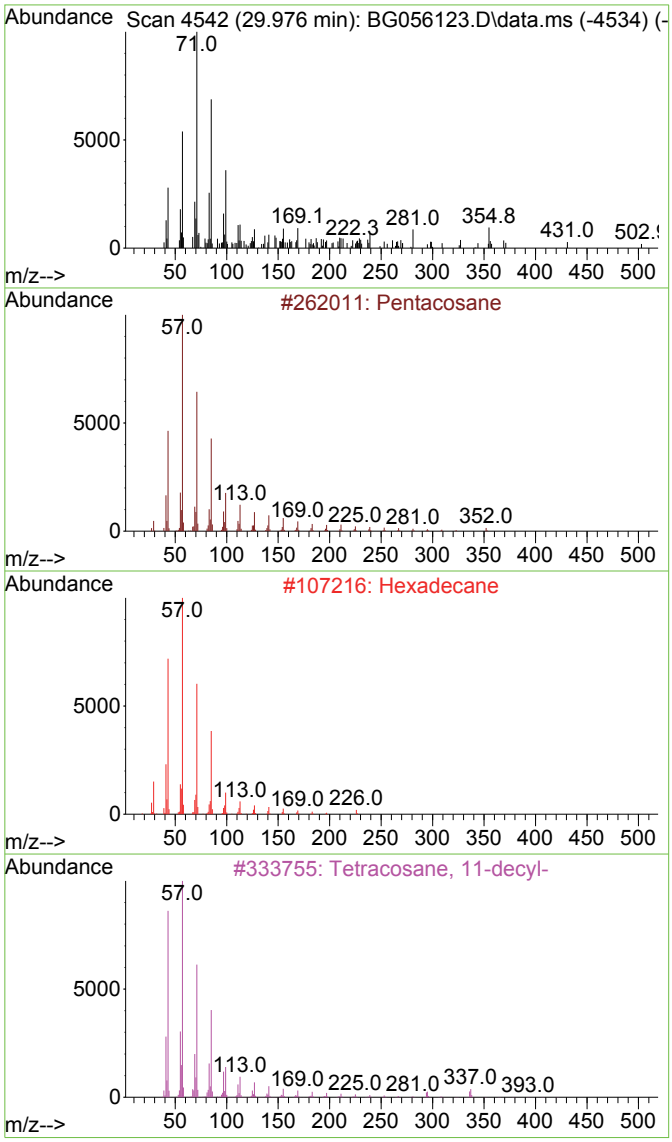
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
Quant Title : SVOA CALIBRATION

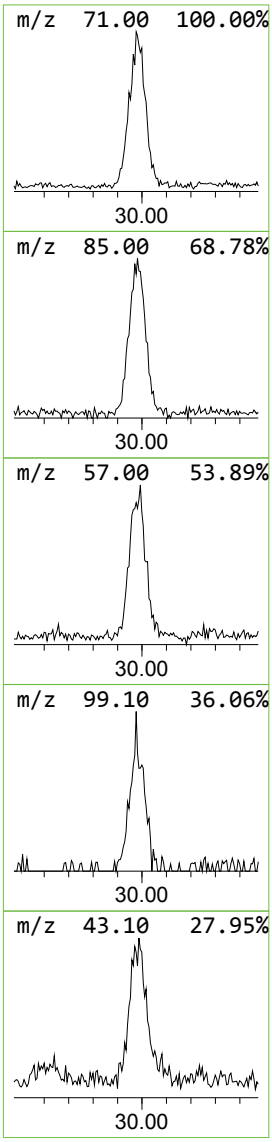
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.976	2.37 ng/ul	102439	Perylene-d12	25.464

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	72
2	Hexadecane	226	C16H34	000544-76-3	68
3	Tetracosane, 11-decyl-	479	C34H70	055429-84-0	64
4	Hexatriacontane	507	C36H74	000630-06-8	64
5	Pentacosane	352	C25H52	000629-99-2	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056123.D
Acq On : 25 Dec 2022 16:12
Operator : CG/JU
Sample : N6017-21
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBYG9

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...	3.372	91.8	ng/ul	963311	1	8.342	209792	20.0
unknown-01	5.381	2.7	ng/ul	28295	1	8.342	209792	20.0
Aceto phenone	9.171	2.1	ng/ul	22225	1	8.342	209792	20.0
Tridecanoic acid	18.536	4.4	ng/ul	157933	4	17.690	720613	20.0
(1H)Pyrrole, 3,...	20.182	5.4	ng/ul	225606	5	21.991	839682	20.0
Heptadecyl hept...	21.574	3.8	ng/ul	157670	5	21.991	839682	20.0
(DEL) Alkane: S...	24.429	2.3	ng/ul	98103	6	25.464	865450	20.0
(DEL) Alkane: S...	26.697	5.0	ng/ul	213987	6	25.464	865450	20.0
Nonacosan-10-ol	26.838	8.3	ng/ul	359189	6	25.464	865450	20.0
(DEL) Alkane: S...	29.976	2.4	ng/ul	102439	6	25.464	865450	20.0