LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
Data File : BG061574.D
Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BHBT1

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-BG052124.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG061574.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.076	8	15	33	rVB	792474	1590366	100.00%	7.868%
2	3.746	123	129	142	rBV	31924	62736	3.94%	0.310%
3	3.852	142	147	153	rVV2	12936	19609	1.23%	0.097%
4	4.992	334	341	350	rVB2	15272	26740	1.68%	0.132%
5	6.555	602	607	614	rVB3	16926	29364	1.85%	0.145%
6	6.860	653	659	664	rBB3	29096	48031	3.02%	0.238%
7	7.072	689	695	705	rVB	113955	202698	12.75%	1.003%
8	7.201	710	717	724	rBV	73003	116907	7.35%	0.578%
9	7.412	747	753	758	rBV	109958	187140	11.77%	0.926%
10	7.459	758	761	769	rVB	15912	27950	1.76%	0.138%
11	7.871	822	831	838	rBV	109369	179242	11.27%	0.887%
12	8.599	948	955	964	rBV	135666	234463	14.74%	1.160%
13	9.028	1021	1028	1036	rBV	110655	198029	12.45%	0.980%
14	9.757	1145	1152	1159	rBV	102524	181115	11.39%	0.896%
15	10.080	1199	1207	1212	rVB6	11397	23458	1.48%	0.116%
16	10.309	1239	1246	1258	rVB	151382	265457	16.69%	1.313%
17	10.667	1297	1307	1313	rBV	142528	250727	15.77%	1.240%
18	10.814	1325	1332	1338	rBV2	28147	58304	3.67%	0.288%
19	12.330	1585	1590	1596	rVB	10960	17146	1.08%	0.085%
20	12.548	1621	1627	1632	rVB2	11101	17253	1.08%	0.085%
21	13.329	1754	1760	1766	rBV3	11648	21444	1.35%	0.106%
22	13.411	1766	1774	1781	rVV	249849	407724	25.64%	2.017%
23	13.817	1832	1843	1850	rBV	137785	242170	15.23%	1.198%
24	13.917	1850	1860	1866	rVV	251163	378513	23.80%	1.873%
25	14.204	1902	1909	1920	rVV	271238	433037	27.23%	2.142%
26	14.469	1948	1954	1957	rBV3	37858	60226	3.79%	0.298%
27	14.510	1957	1961	1967	rVV	222327	339822	21.37%	1.681%
28	14.575	1967	1972	1980	rVB5	17203	34001	2.14%	0.168%
29	14.727	1991	1998	2002	rBV3	34515	84221	5.30%	0.417%
30	14.774	2002	2006	2012	rVB	67592	112091	7.05%	0.555%
31	14.915	2020	2030	2038	rVB2	20795	53430	3.36%	0.264%
32	14.992	2038	2043	2046	rVV3	15154	21418	1.35%	0.106%
33	15.127	2056	2066	2072	rBV6	10607	25011	1.57%	0.124%
34	15.303	2090	2096	2100	rBV5	10559	21425	1.35%	0.106%
35	15.509	2122	2131	2136	rBV	355832	541327	34.04%	2.678%
36	15.562	2136	2140	2145	rVB2	21858	31555	1.98%	0.156%

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Misc :
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Instrument :
BNA_G
ClientSampleId :
BHBT1

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-BG052124.MA.M
Title : SVOA CALIBRATION

37	15.662	2151	2157	2164	rBV	73665	120433	7.57%	0.596%
38	15.855	2185	2190	2199	rVV7	17087	36696	2.31%	0.182%
39	16.002	2208	2215	2223	rVB7	11672	29276	1.84%	0.145%
40	16.243	2245	2256	2260	rBV6	26888	57372	3.61%	0.284%
41	16.766	2341	2345	2348	rBV4	12149	19593	1.23%	0.097%
42	16.925	2366	2372	2378	rBV2	28071	45358	2.85%	0.224%
43	17.013	2380	2387	2392	rBV6	28127	49250	3.10%	0.244%
44	17.083	2392	2399	2408	rVB2	21743	43629	2.74%	0.216%
45	17.266	2419	2430	2434	rBV	302265	475116	29.87%	2.351%
46	17.307	2434	2437	2442	rVB	328796	448132	28.18%	2.217%
47	17.365	2442	2447	2451	rBV	395534	561532	35.31%	2.778%
48	17.448	2457	2461	2467	rBV5	12510	25187	1.58%	0.125%
49	17.583	2481	2484	2489	rVB5	12736	21313	1.34%	0.105%
50	17.677	2491	2500	2504	rBV2	26454	53261	3.35%	0.263%
51	17.759	2510	2514	2516	rBV4	12892	17579	1.11%	0.087%
52	17.847	2525	2529	2536	rVB3	21071	39677	2.49%	0.196%
53	17.929	2539	2543	2545	rBV5	14720	17825	1.12%	0.088%
54	18.164	2574	2583	2585	rBV2	150483	312919	19.68%	1.548%
55	18.194	2585	2588	2592	rVB	117002	169895	10.68%	0.841%
56	18.276	2597	2602	2607	rVV	24326	50315	3.16%	0.249%
57	18.335	2607	2612	2616	rVV2	101971	201133	12.65%	0.995%
58	18.376	2616	2619	2625	rVB	65771	95044	5.98%	0.470%
59	18.452	2627	2632	2636	rBV4	21169	37836	2.38%	0.187%
60	18.494	2637	2639	2642	rVV3	14636	18051	1.14%	0.089%
61	18.541	2643	2647	2651	rVB5	16271	25333	1.59%	0.125%
62	18.640	2659	2664	2669	rBV	56427	86766	5.46%	0.429%
63	18.693	2669	2673	2678	rVB2	65141	100056	6.29%	0.495%
64	18.893	2703	2707	2711	rVB2	25122	34962	2.20%	0.173%
65	18.940	2711	2715	2719	rVV	30950	42133	2.65%	0.208%
66	18.981	2719	2722	2726	rVB2	23669	28582	1.80%	0.141%
67	19.087	2731	2740	2749	rBV3	138944	429298	26.99%	2.124%
68	19.157	2749	2752	2755	rVB	27543	30557	1.92%	0.151%
69	19.210	2755	2761	2768	rBV2	61075	132784	8.35%	0.657%
70	19.328	2773	2781	2787	rBV	724110	1046735	65.82%	5.178%
71	19.387	2788	2791	2794	rBV	23633	29965	1.88%	0.148%
72	19.434	2795	2799	2804	rBV6	27166	49611	3.12%	0.245%
73	19.663	2832	2838	2840	rBV2	556448	808369	50.83%	3.999%
74	19.692	2840	2843	2850	rVV	563526	785785	49.41%	3.887%
75	19.763	2851	2855	2862	rVB3	53365	89524	5.63%	0.443%
76	19.933	2880	2884	2891	rVB2	36433	63749	4.01%	0.315%

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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

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

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Title : SVOA CALIBRATION

77	20.015	2892	2898	2904	rBV4	68505	188392	11.85%	0.932%
78	20.150	2916	2921	2922	rBV3	54767	83844	5.27%	0.415%
79	20.186	2924	2927	2932	rVB2	119166	164886	10.37%	0.816%
80	20.286	2940	2944	2948	rBV	75093	108100	6.80%	0.535%
81	20.485	2974	2978	2981	rBV	57177	77492	4.87%	0.383%
82	20.526	2982	2985	2989	rVV2	70996	96277	6.05%	0.476%
83	20.573	2991	2993	2997	rVB	56876	62457	3.93%	0.309%
84	20.644	3002	3005	3008	rVB3	74864	78692	4.95%	0.389%
85	20.761	3023	3025	3032	rBV8	47853	85108	5.35%	0.421%
86	20.944	3052	3056	3062	rVB	98789	144566	9.09%	0.715%
87	21.002	3062	3066	3070	rBV	205345	263456	16.57%	1.303%
88	21.173	3093	3095	3099	rVB	65716	69926	4.40%	0.346%
89	21.273	3108	3112	3119	rVB	94862	154305	9.70%	0.763%
90	21.408	3131	3135	3140	rVB	193391	269864	16.97%	1.335%
91	21.519	3147	3154	3159	rBV4	469036	1167900	73.44%	5.778%
92	21.572	3159	3163	3176	rVB	367274	625759	39.35%	3.096%
93	22.295	3282	3286	3296	rVB5	77316	231694	14.57%	1.146%
94	23.288	3450	3455	3462	rVB5	133747	258598	16.26%	1.279%
95	23.652	3511	3517	3523	rBV	239641	678105	42.64%	3.355%
96	23.717	3525	3528	3534	rVV3	242036	435702	27.40%	2.156%
97	23.793	3535	3541	3552	rVB4	129760	372415	23.42%	1.842%
98	24.422	3642	3648	3655	rBV	149891	362144	22.77%	1.792%
99	24.627	3677	3683	3694	rVB	169080	442276	27.81%	2.188%
100	25.691	3857	3864	3874	rVB	183081	516656	32.49%	2.556%

Sum of corrected areas: 20213395

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
Data File : BG061574.D
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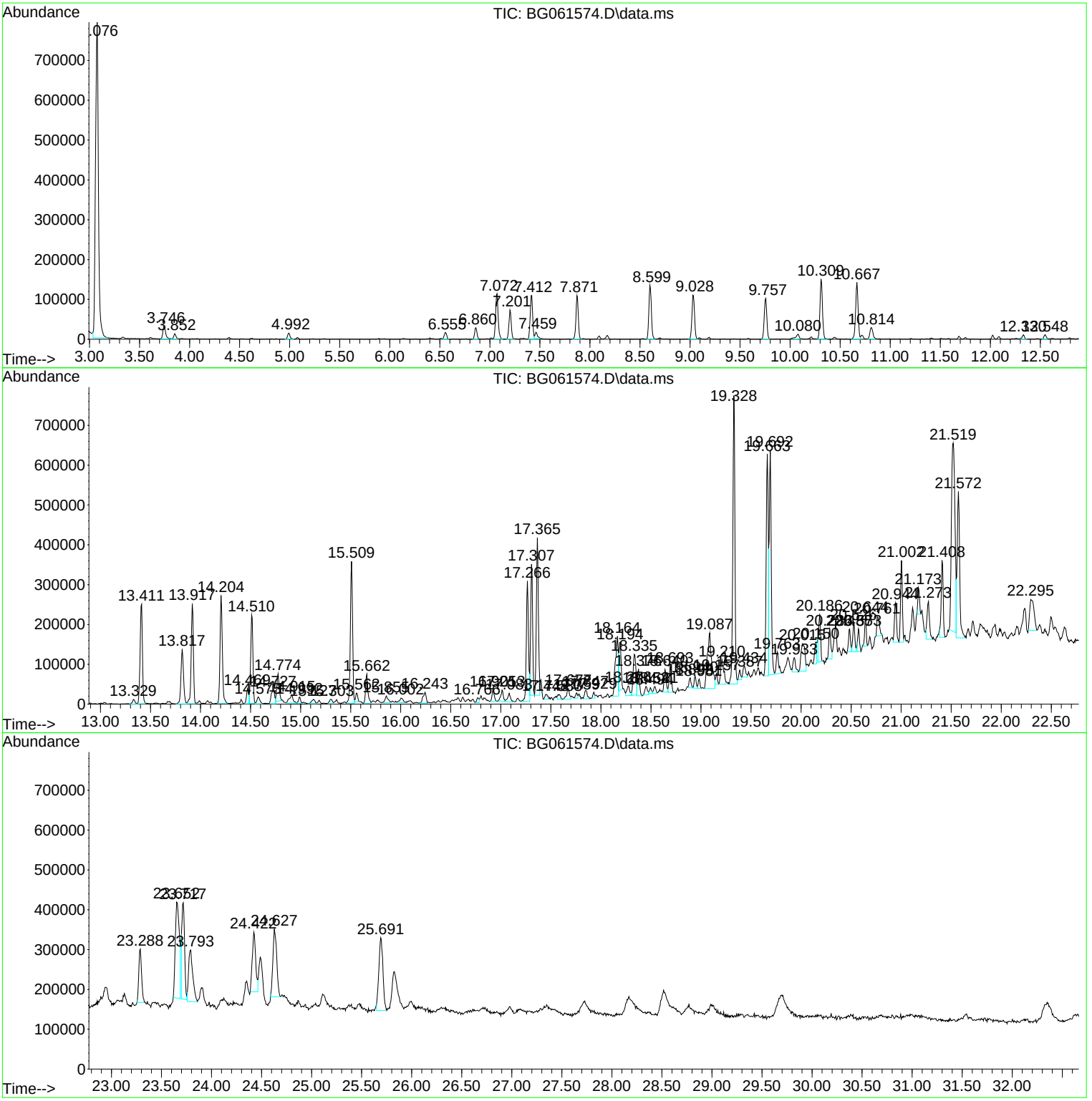
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
Data File : BG061574.D
Acq On : 8 Jun 2024 21:00
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Sample : P2647-17
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

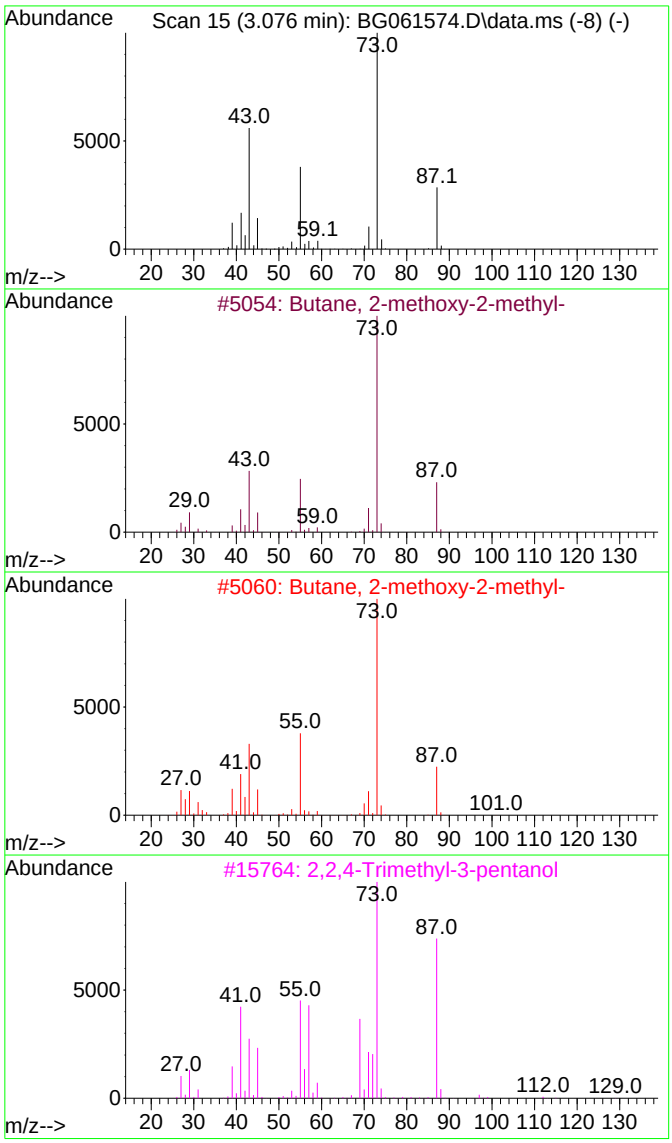
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

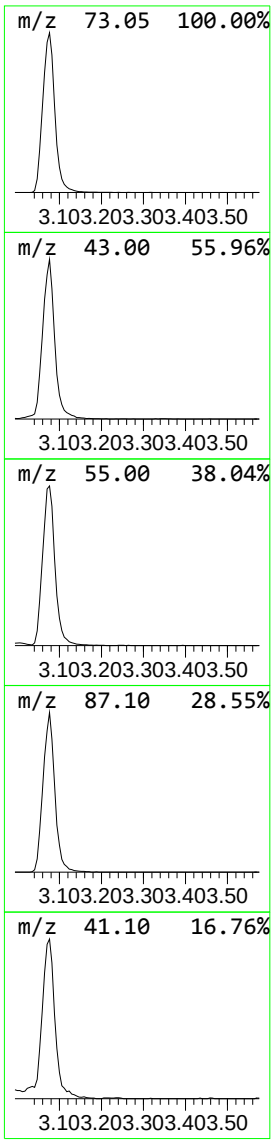
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.076	177.46 ng/ul	1590370	1,4-Dichlorobenzene-d4	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	36
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33
5	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
Data File : BG061574.D
Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

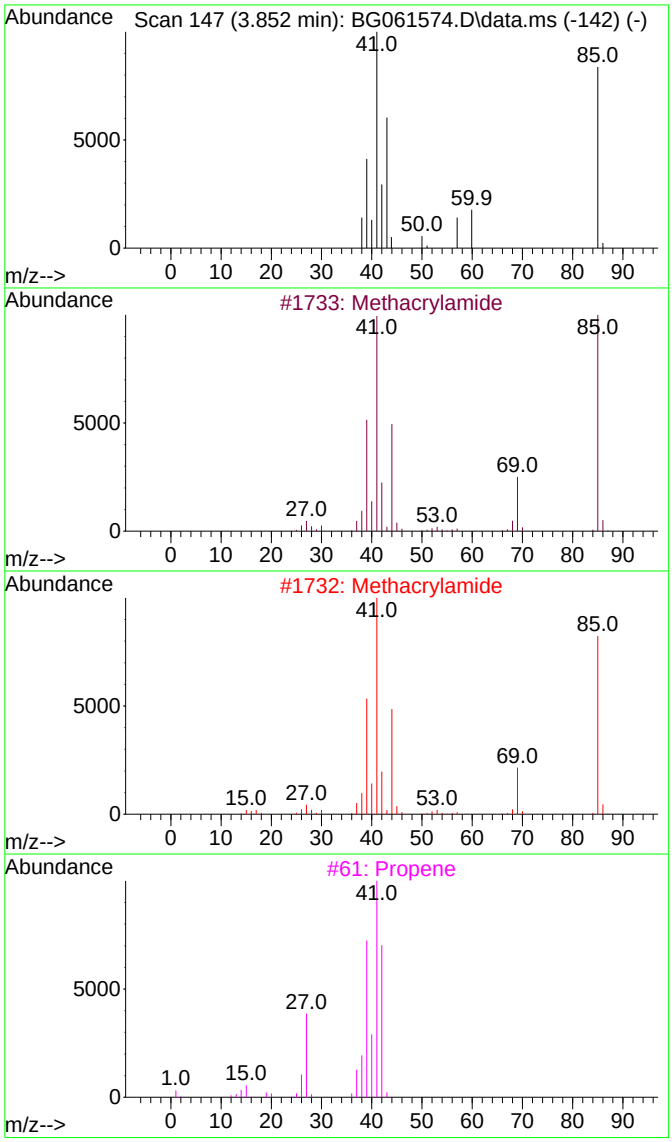
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

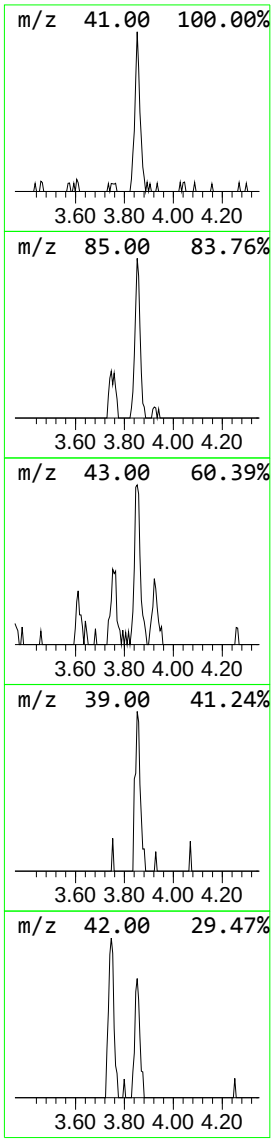
****

Peak Number 2 unknown-01 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.852	2.19 ng/ul	19609	1,4-Dichlorobenzene-d4	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methacrylamide	85	C4H7NO	000079-39-0	39
2	Methacrylamide	85	C4H7NO	000079-39-0	33
3	Propene	42	C3H6	000115-07-1	9
4	trans-Crotonamide	85	C4H7NO	000625-37-6	7
5	Propene	42	C3H6	000115-07-1	7





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
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Instrument :
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ClientSampleId :
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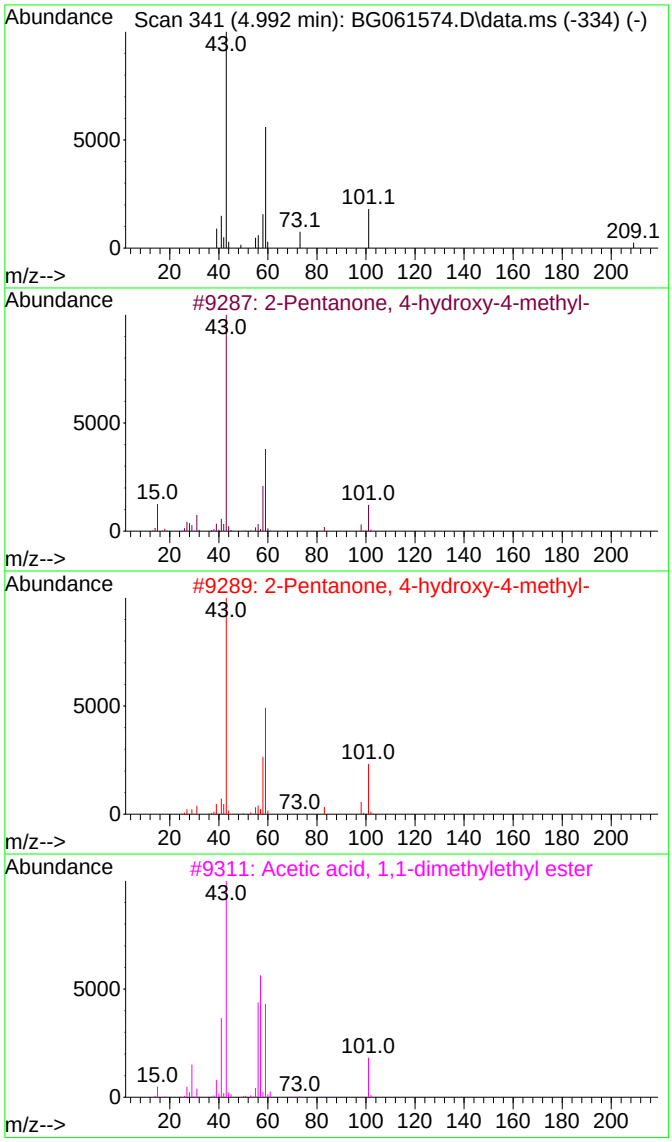
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

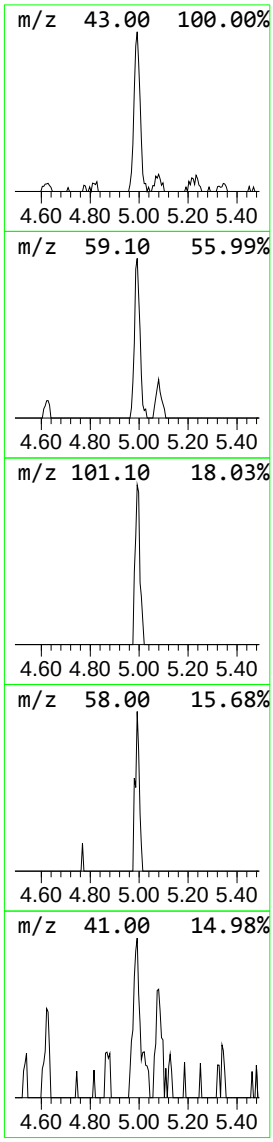
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.992	2.98 ng/ul	26740	1,4-Dichlorobenzene-d4	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4	3-Octanol	130	C8H18O	000589-98-0	36
5	Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

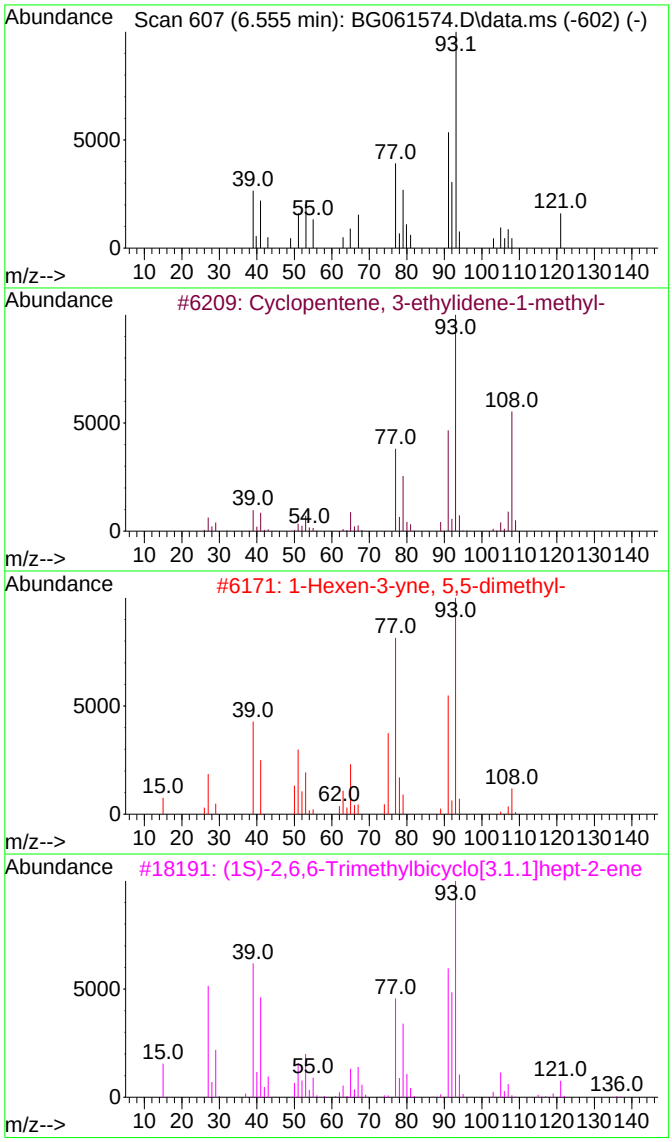
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

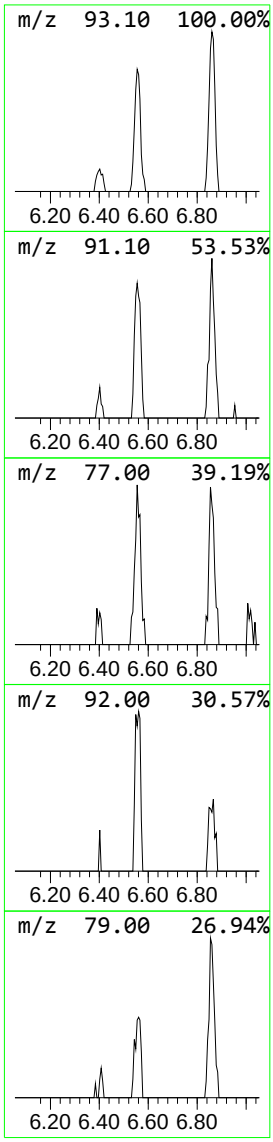
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclopentene, 3-ethylidene-... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.555	3.28 ng/ul	29364	1,4-Dichlorobenzene-d4	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 3-ethylidene-1-met...	108	C8H12	062338-00-5	53
2	1-Hexen-3-yne, 5,5-dimethyl-	108	C8H12	004911-58-4	50
3	(1S)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-26-4	50
4	trans-.beta.-Ocimene	136	C10H16	003779-61-1	47
5	1,4-Methano-1H-Cyclopropa[d]pyri...	136	C8H12N2	109746-10-3	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
Data File : BG061574.D
Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

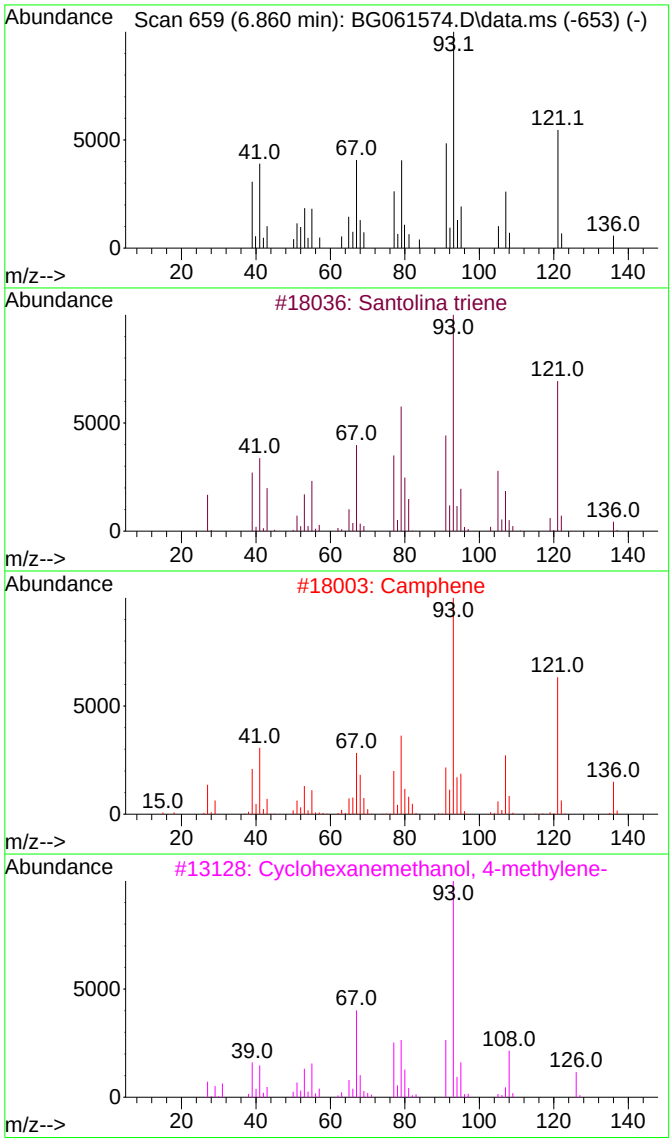
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

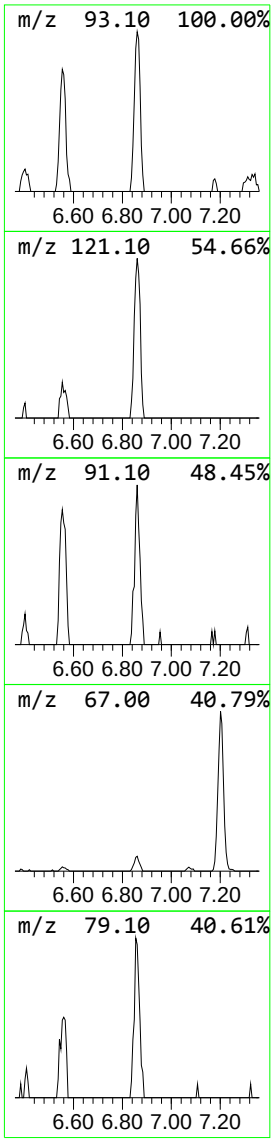
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Santolina triene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.860	5.36 ng/ul	48031	1,4-Dichlorobenzene-d4	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Santolina triene	136	C10H16	002153-66-4	93
2	Camphene	136	C10H16	000079-92-5	72
3	Cyclohexanemethanol, 4-methylene-	126	C8H14O	001004-24-6	59
4	.alpha.-Pinene	136	C10H16	000080-56-8	53
5	1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	53





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

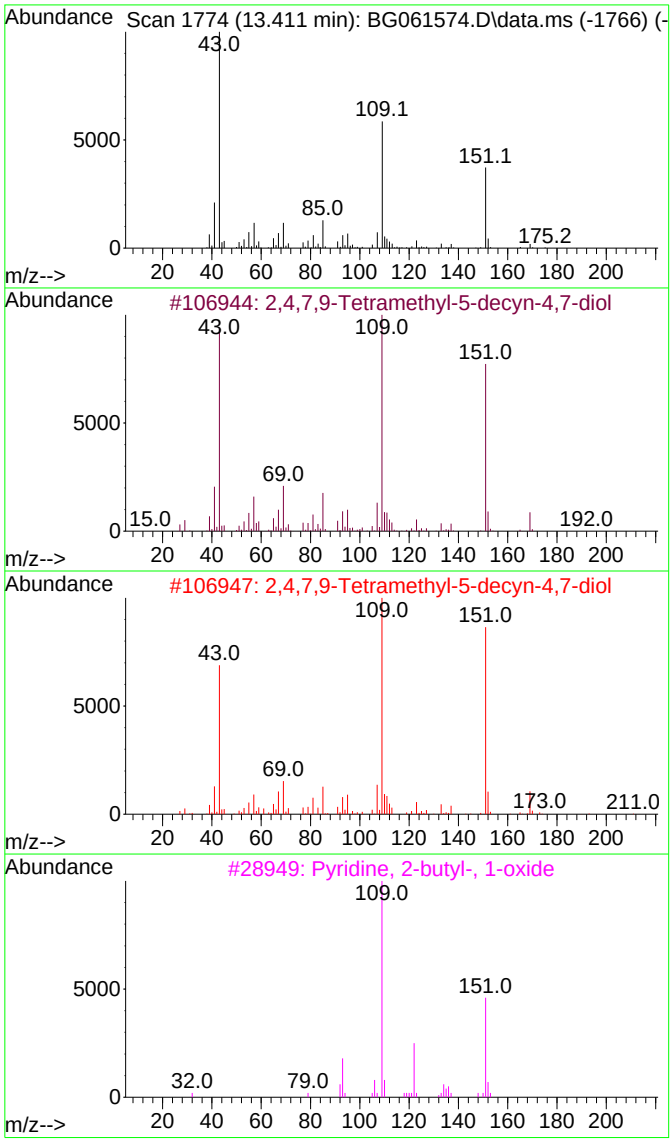
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

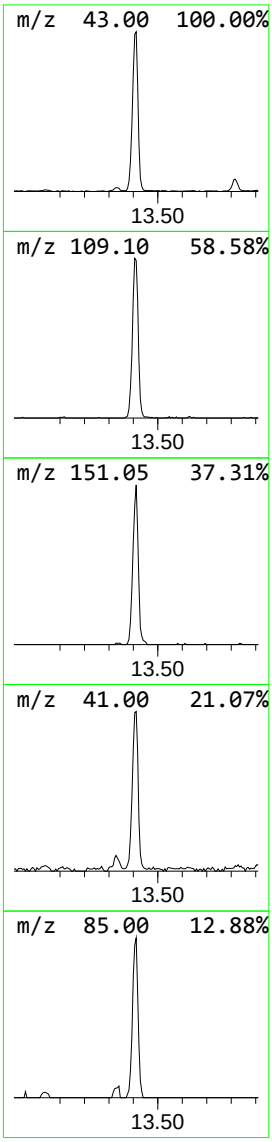
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.411	24.00 ng/ul	407724	Acenaphthene-d10	14.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	87
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	83
3	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	53
4	Acetaminophen	151	C8H9NO2	000103-90-2	47
5	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	47





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

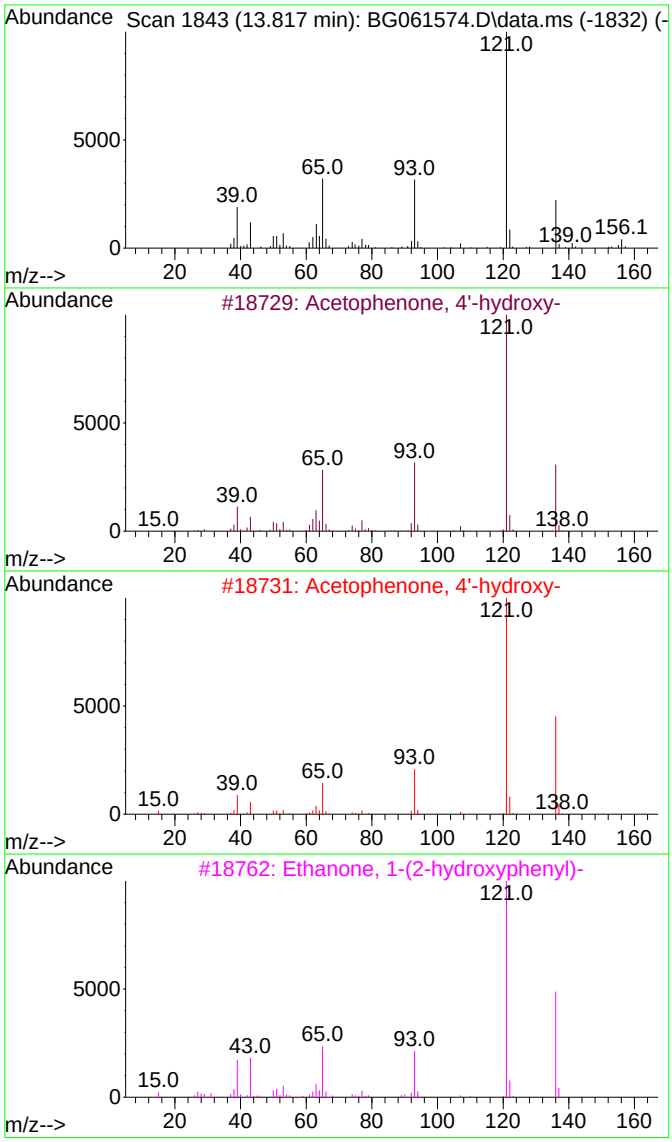
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

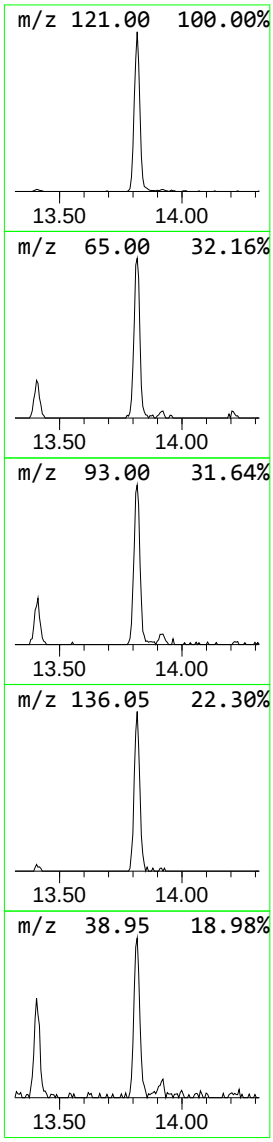
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Acetophenone, 4'-hydroxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.817	14.25 ng/ul	242170	Acenaphthene-d10	14.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone, 4'-hydroxy-	136	C8H8O2	000099-93-4	96
2	Acetophenone, 4'-hydroxy-	136	C8H8O2	000099-93-4	95
3	Ethanone, 1-(2-hydroxyphenyl)-	136	C8H8O2	000118-93-4	94
4	Acetophenone, 4'-hydroxy-	136	C8H8O2	000099-93-4	94
5	Ethanone, 1-(2-hydroxyphenyl)-	136	C8H8O2	000118-93-4	91





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

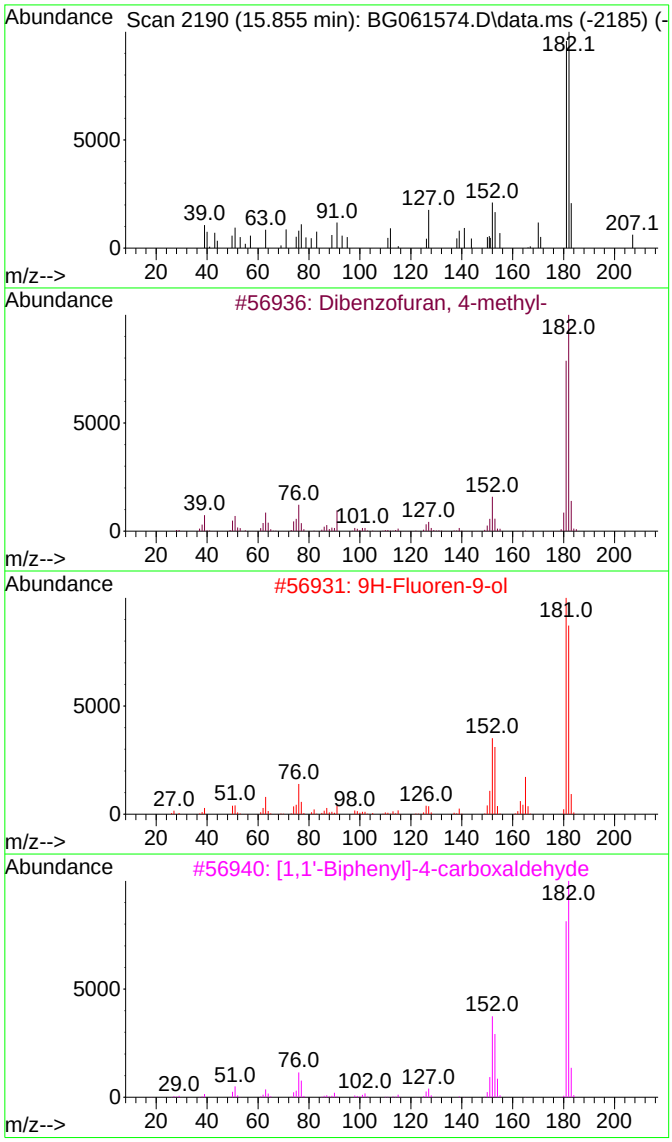
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

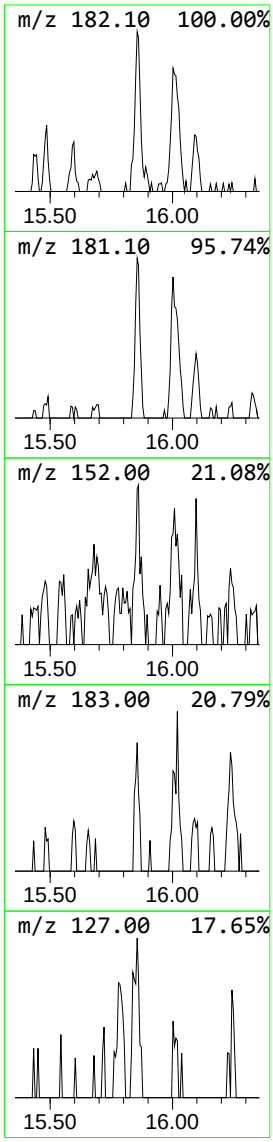
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Dibenzofuran, 4-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.855	2.16 ng/ul	36696	Acenaphthene-d10	14.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	80
2	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64
3	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	64
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	64
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	64





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Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

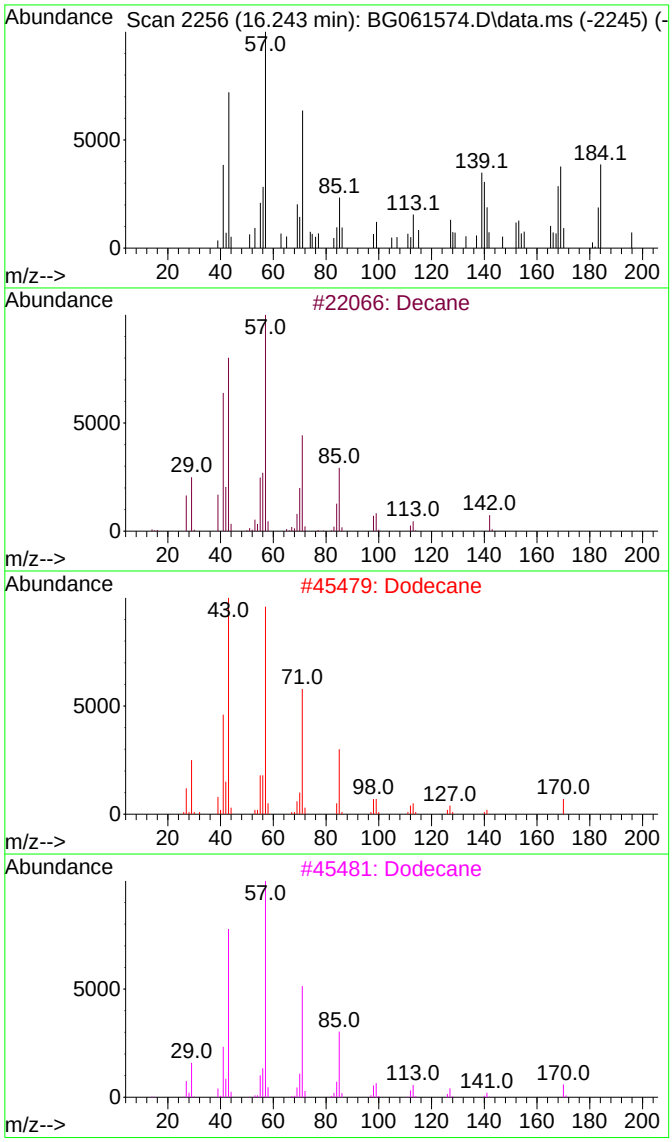
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

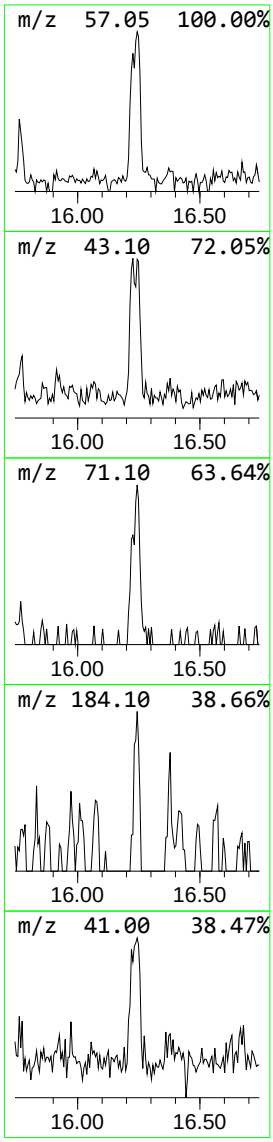
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.243	2.42 ng/ul	57372	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	25
2	Dodecane	170	C12H26	000112-40-3	22
3	Dodecane	170	C12H26	000112-40-3	15
4	Decane	142	C10H22	000124-18-5	15
5	Oxalic acid, isohexyl neopentyl ...	244	C13H24O4	1000309-73-0	14





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

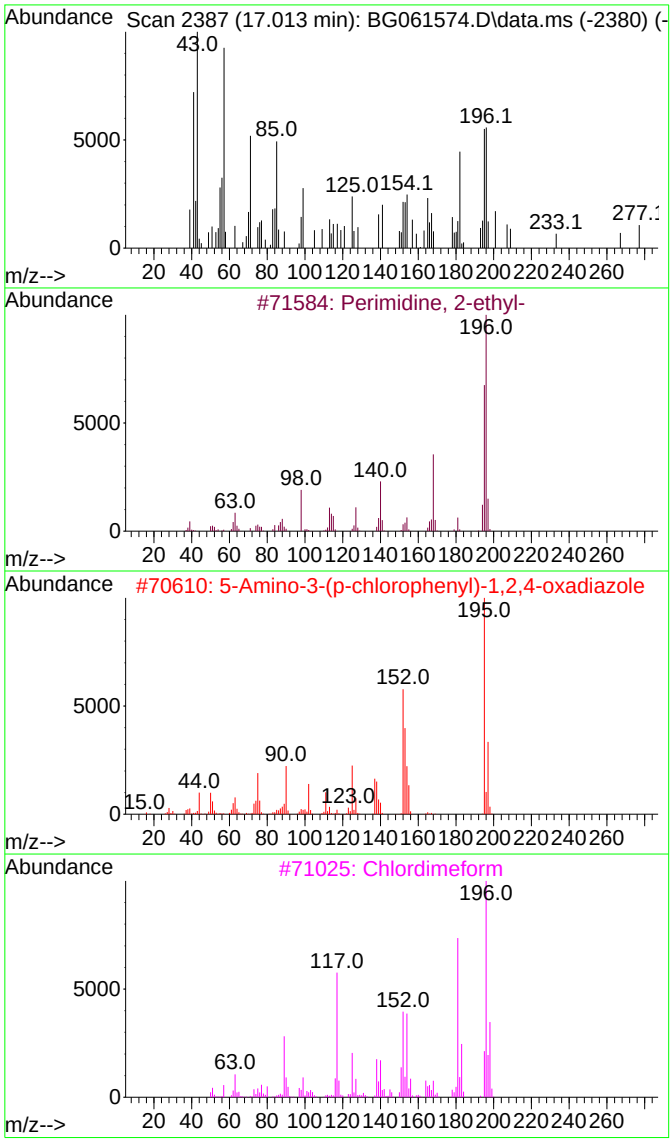
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

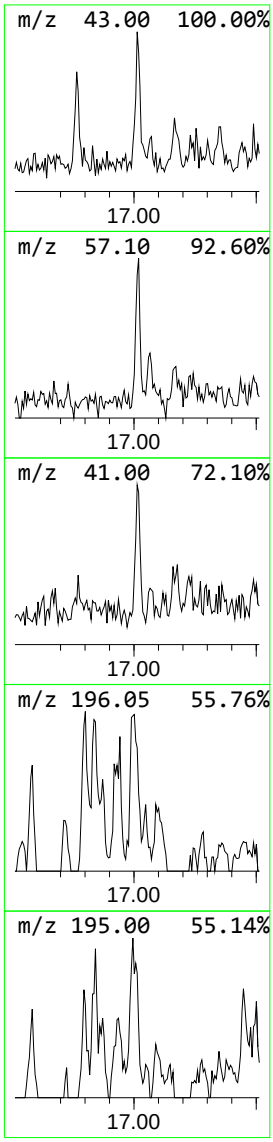
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 unknown-02 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.013	2.07 ng/ul	49250	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Perimidine, 2-ethyl-	196	C13H12N2	028478-13-9	25
2	5-Amino-3-(p-chlorophenyl)-1,2,4...	195	C8H6ClN3O	114212-86-1	22
3	Chlordimeform	196	C10H13ClN2	006164-98-3	22
4	Benzo[b]benzofuran-2-carboxaldehyde	196	C13H8O2	1000351-56-0	18
5	N-[(7-Methoxy-2H-1,3-benzodioxol...	195	C9H9NO4	1000387-37-6	18





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

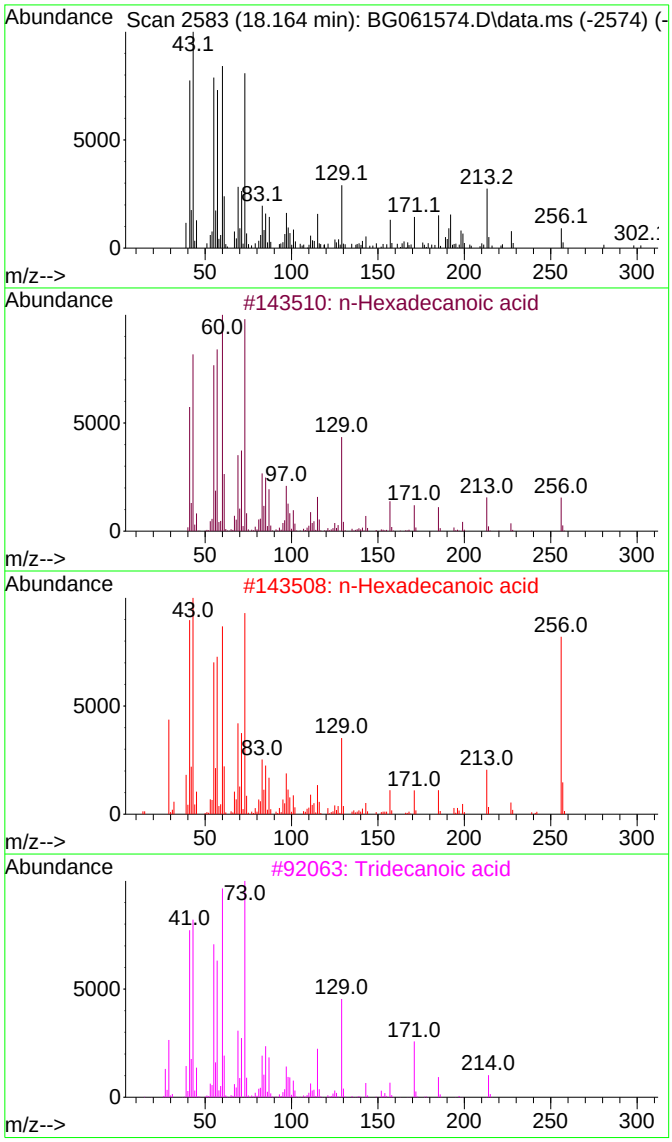
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

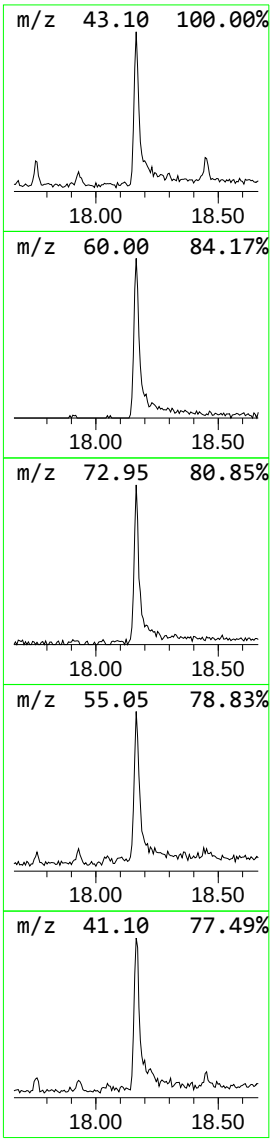
****

Peak Number 11 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.165	13.17 ng/ul	312919	Phenanthrene-d10	17.265

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	97
3	Tridecanoic acid	214	C13H26O2	000638-53-9	93
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
5	Nonanoic acid	158	C9H18O2	000112-05-0	64





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

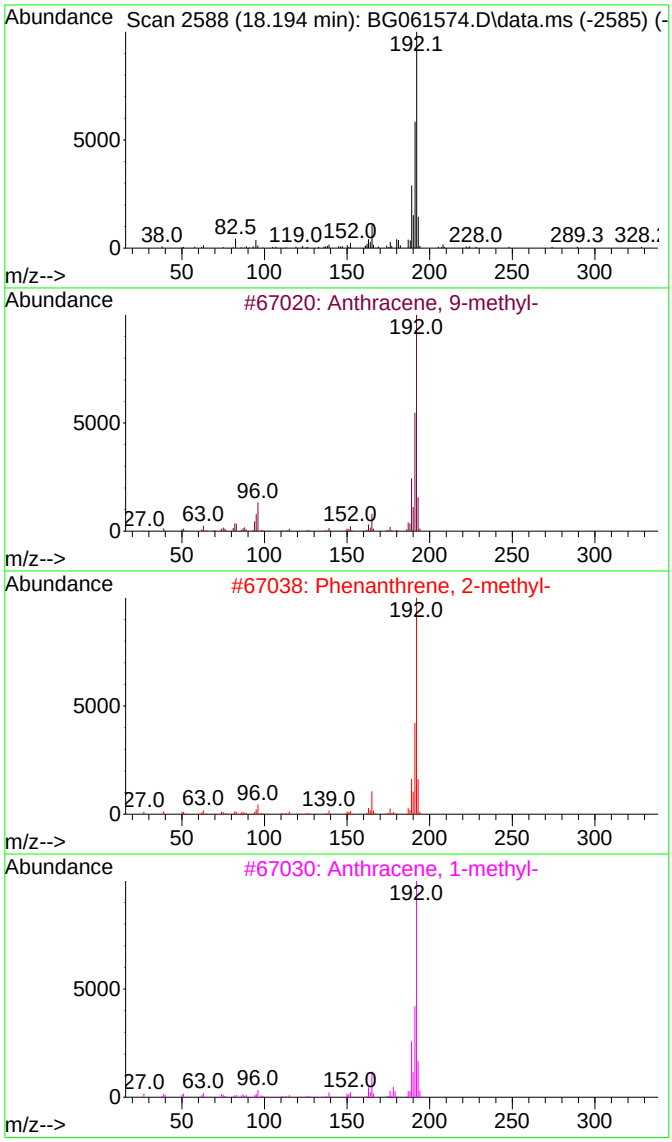
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

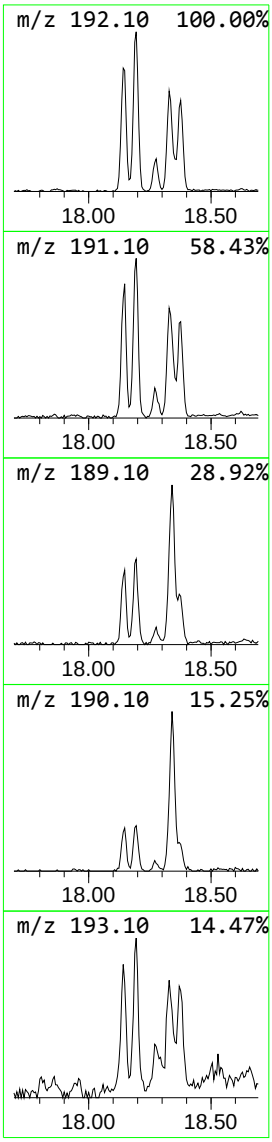
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Anthracene, 9-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.194	7.15 ng/ul	169895	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

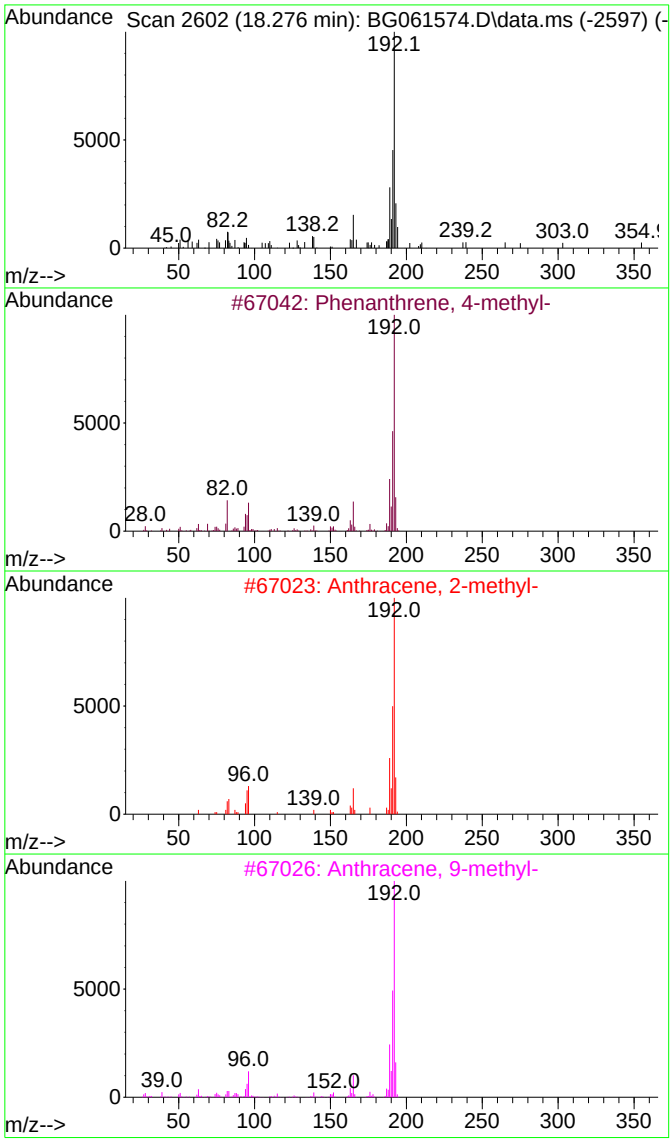
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

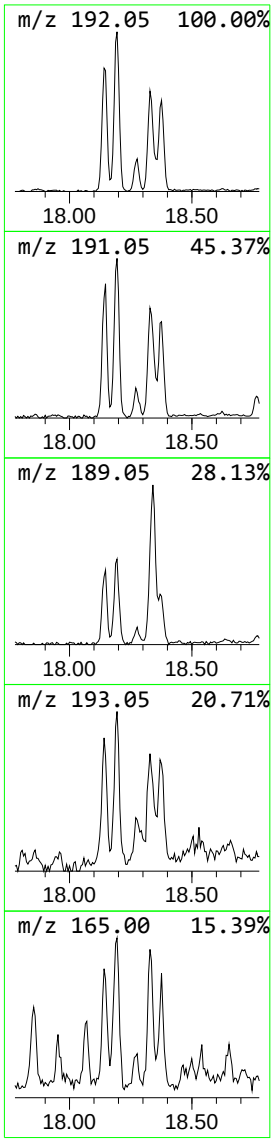
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phenanthrene, 4-methyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.276	2.12 ng/ul	50315	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	76
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	74
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	74
4	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	72
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	72





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

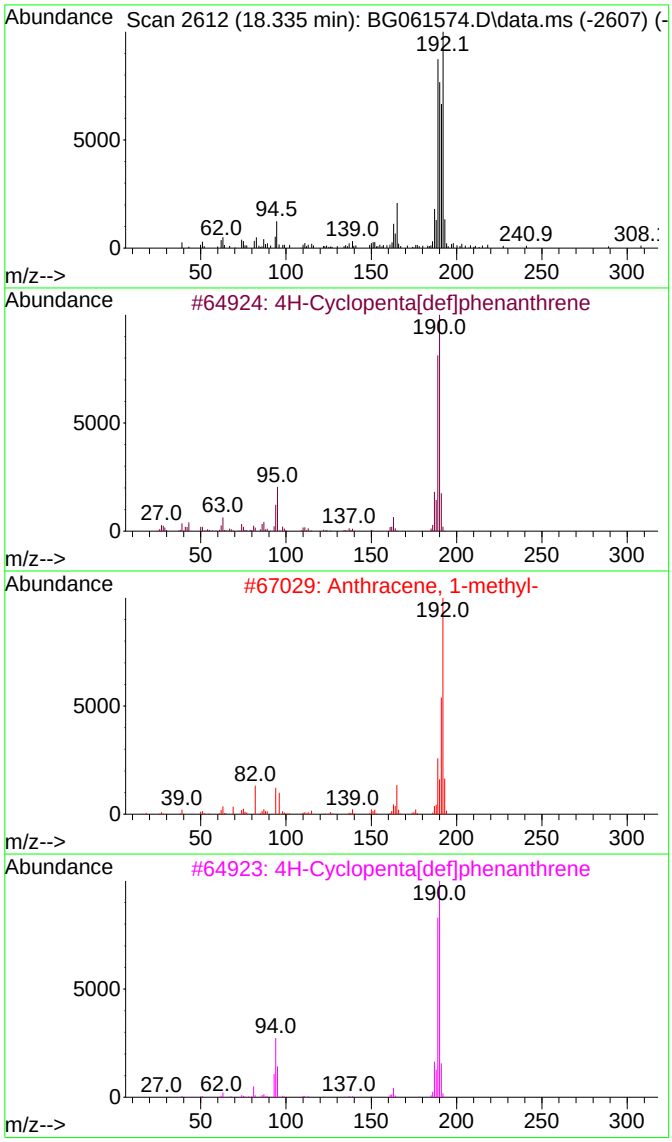
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

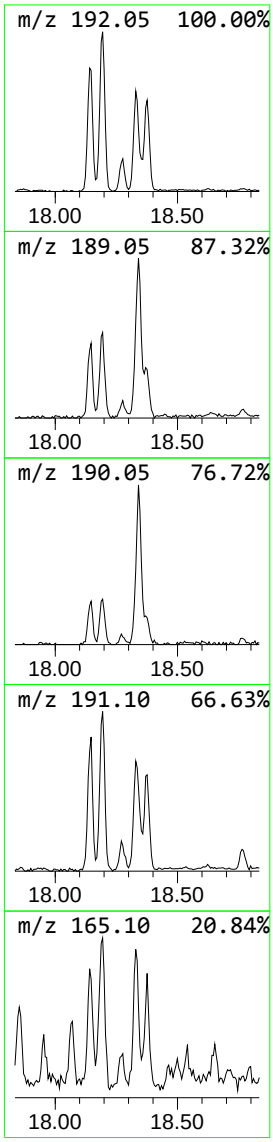
****

Peak Number 14 unknown-03 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.335	8.47 ng/ul	201133	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	45
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	41
5	Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	38





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Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

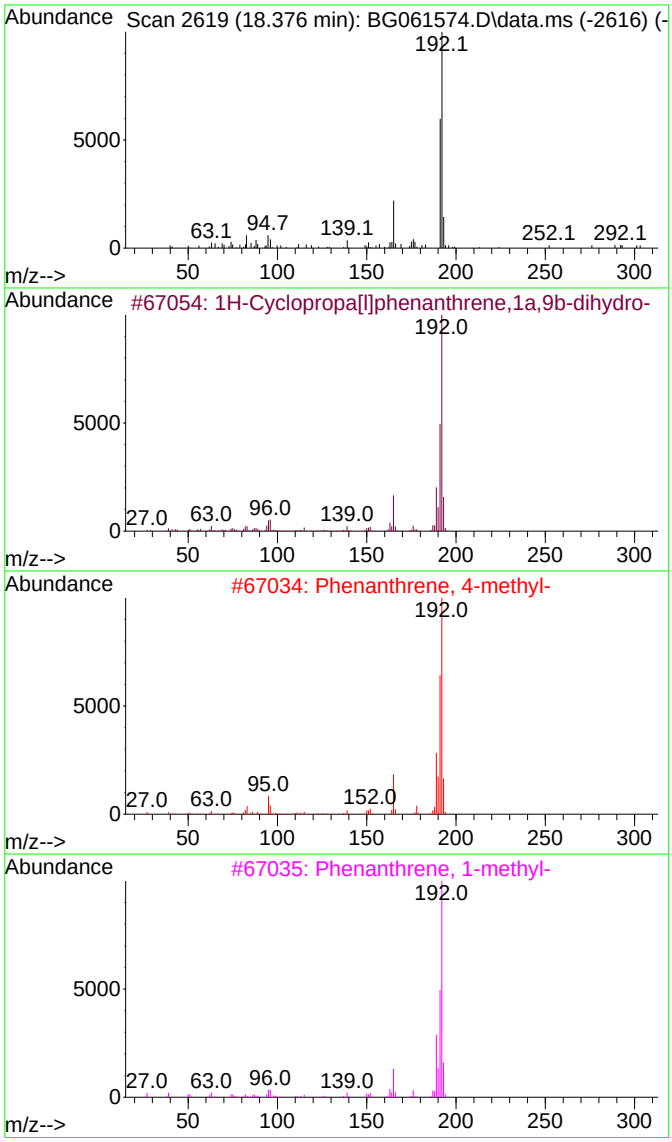
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

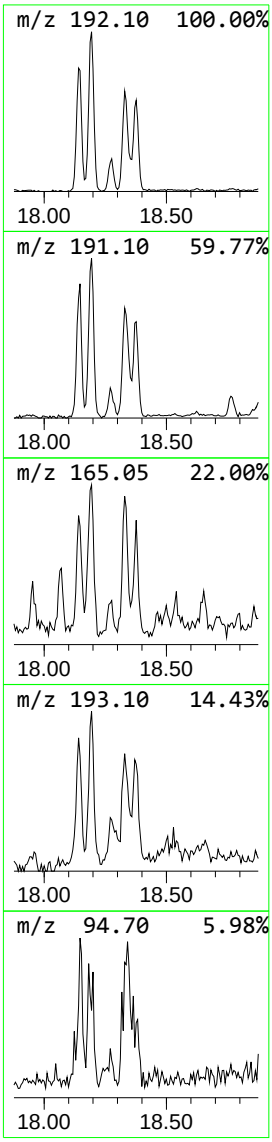
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 1H-Cyclopropa[l]phenanthren... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.376	4.00 ng/ul	95044	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	86
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	86





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

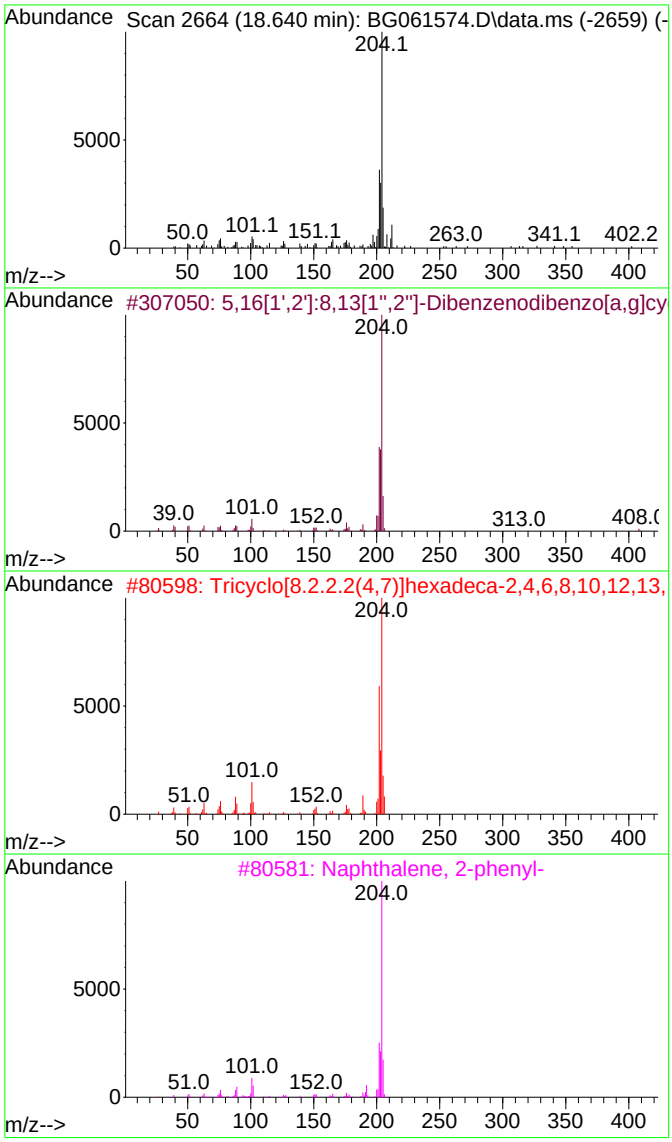
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

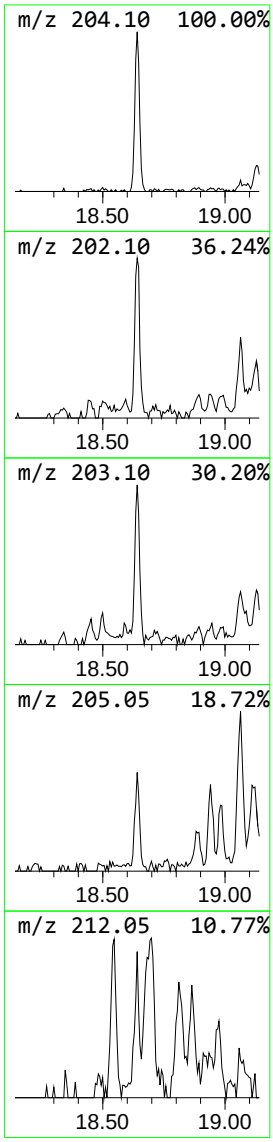
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.640	3.65 ng/ul	86766	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	74
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	70
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	62
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	62





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Data File : BG061574.D
Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

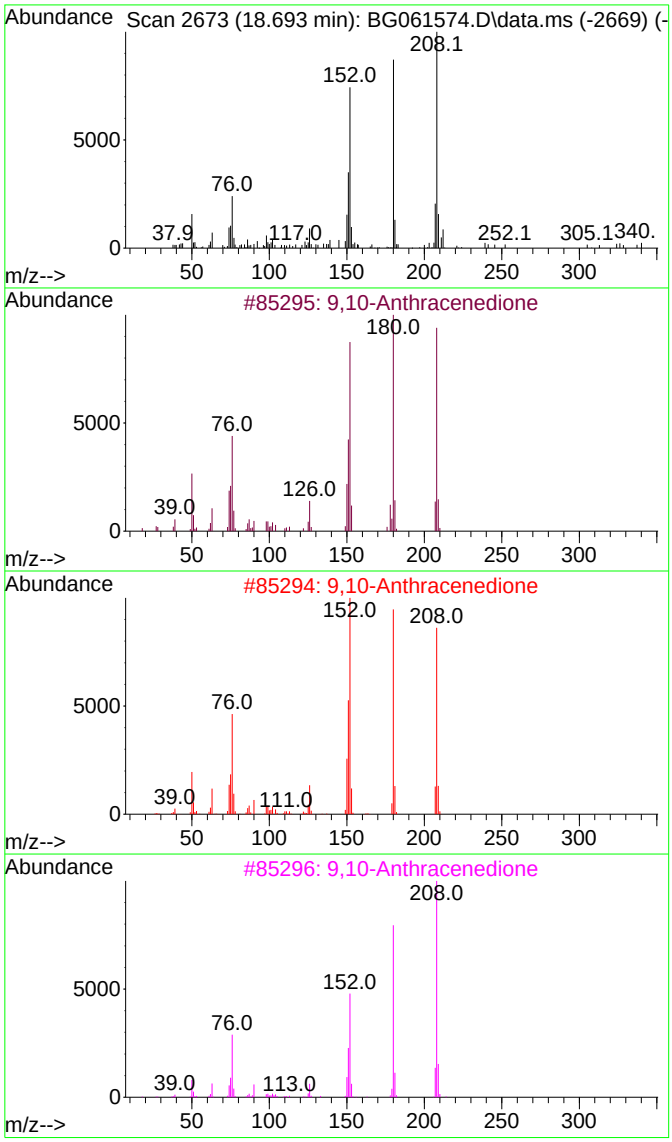
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

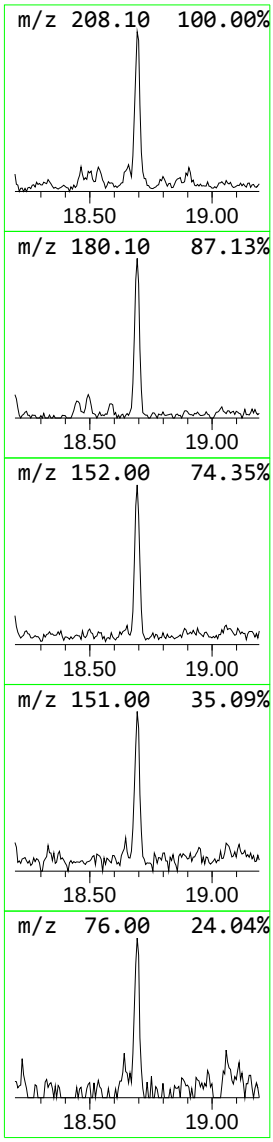
****

Peak Number 17 9,10-Anthracenedione Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.693	4.21 ng/ul	100056	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
5	1H-Phenalen-1-one	180	C13H8O	000548-39-0	78





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Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

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Quant Title : SVOA CALIBRATION

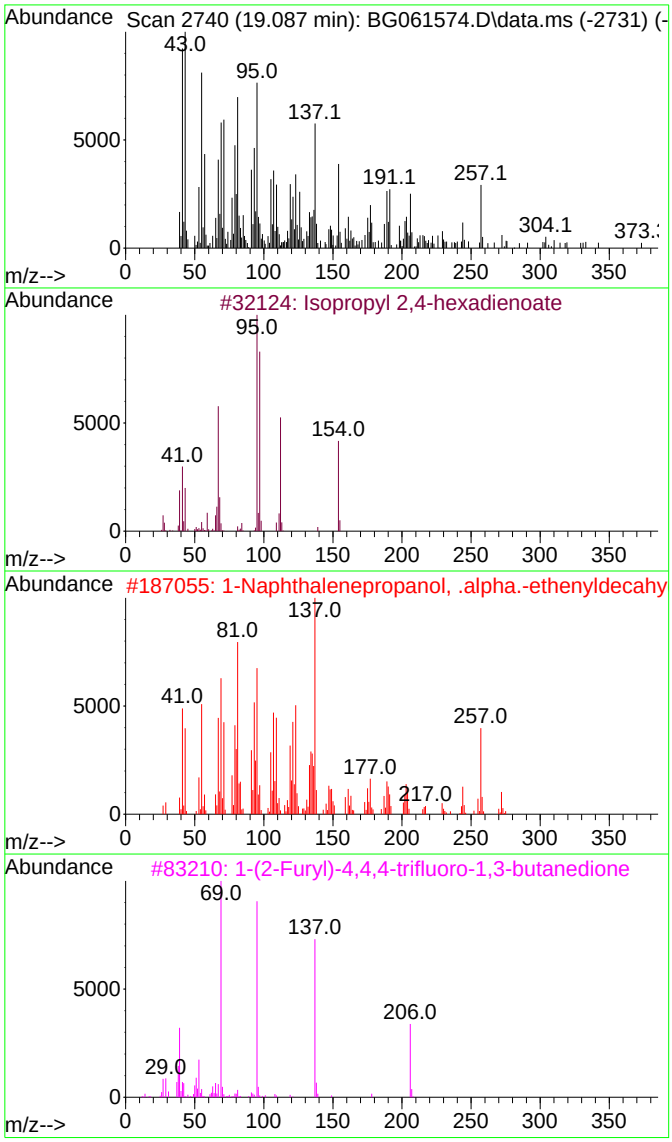
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

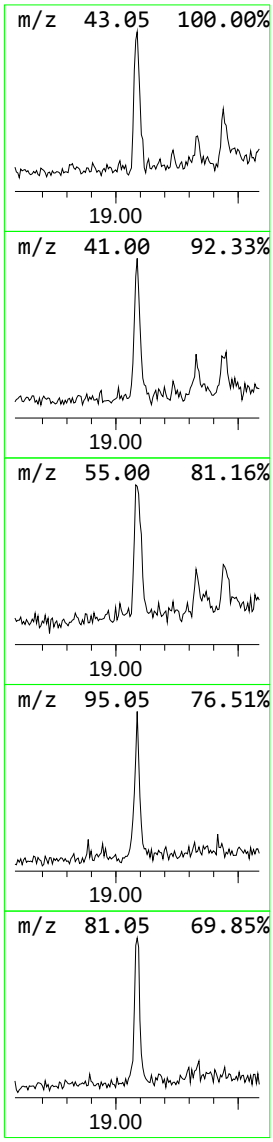
****

Peak Number 18 unknown-04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.087	18.07 ng/ul	429298	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl 2,4-hexadienoate	154	C9H14O2	044987-75-9	42
2	1-Naphthalenepropanol, .alpha.-e...	290	C20H34O	001438-62-6	42
3	1-(2-Furyl)-4,4,4-trifluoro-1,3-...	206	C8H5F3O3	000326-90-9	38
4	1-Naphthalenepropanol, .alpha.-e...	290	C20H34O	001438-62-6	38
5	cis, cis-3-Ethylbicyclo[4.4.0]de...	166	C12H22	066660-42-2	38





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Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

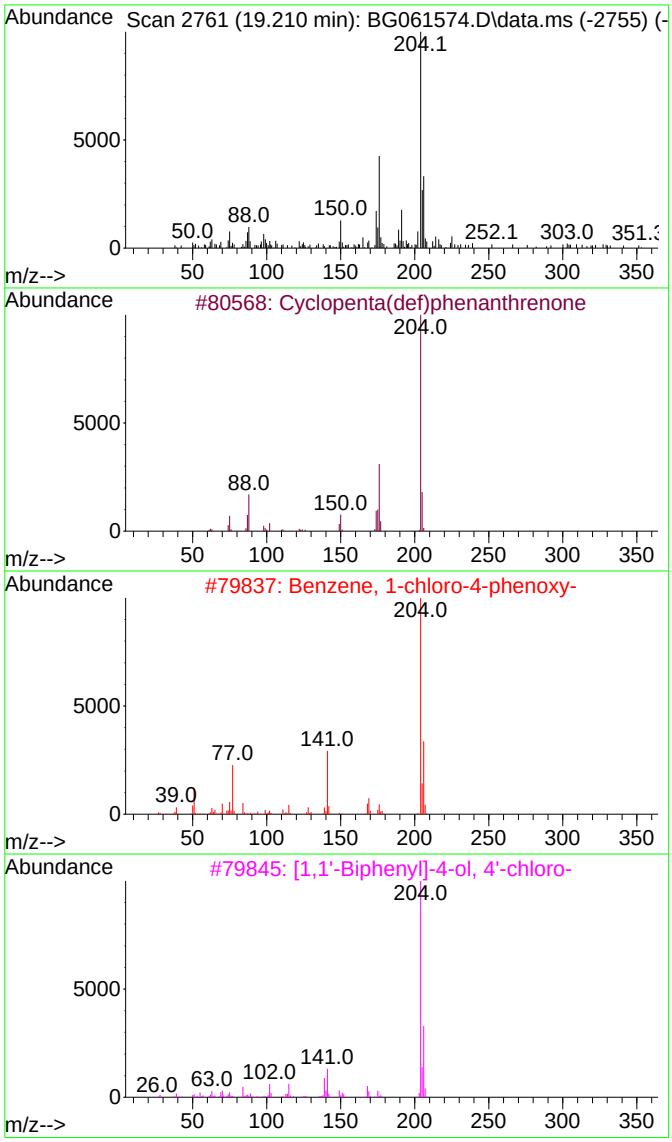
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

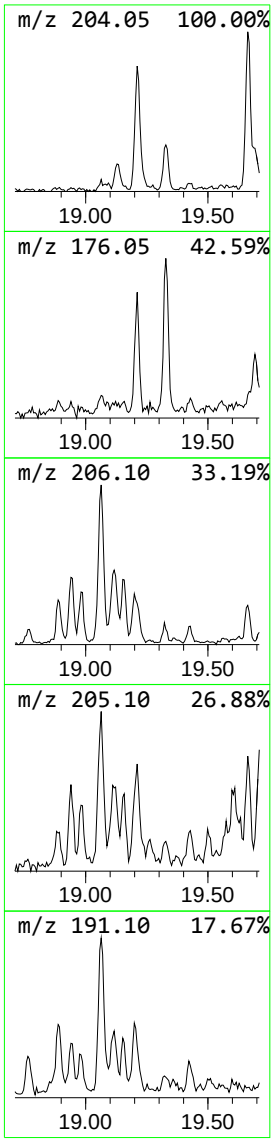
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Cyclopenta(def)phenanthrenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.210	5.59 ng/ul	132784	Phenanthrene-d10	17.265

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	52
2	Benzene, 1-chloro-4-phenoxy-	204	C12H9ClO	007005-72-3	46
3	[1,1'-Biphenyl]-4-ol, 4'-chloro-	204	C12H9ClO	028034-99-3	43
4	[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	43
5	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	43





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Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

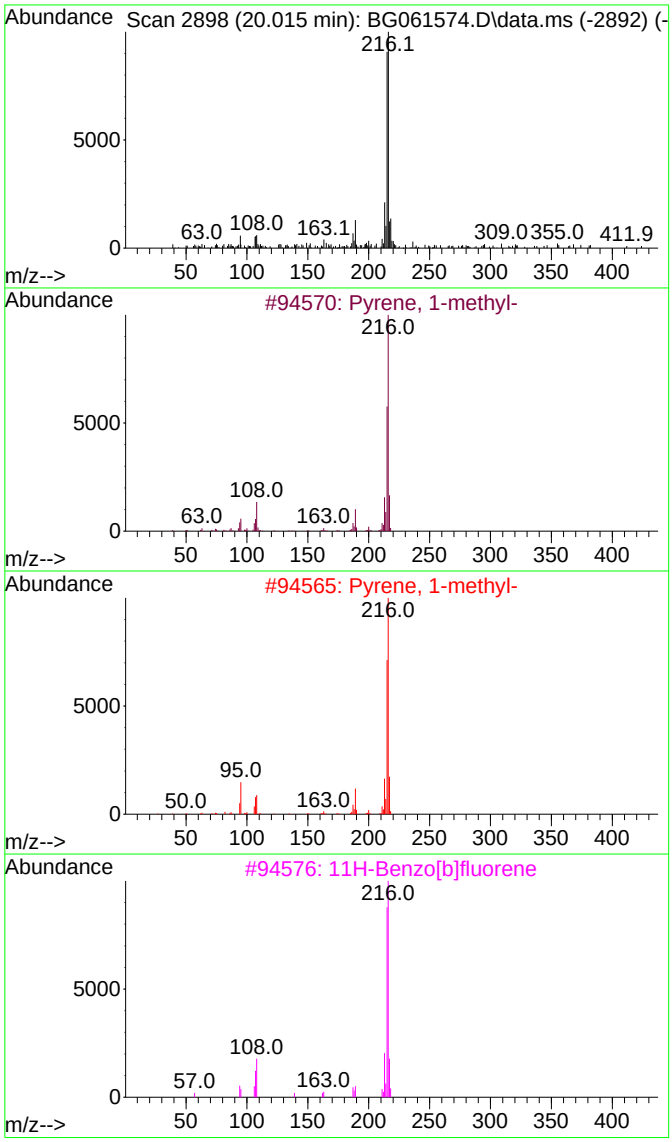
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

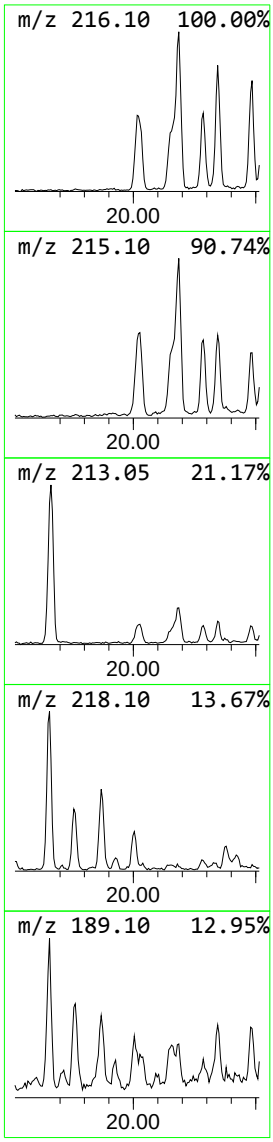
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Pyrene, 1-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.015	3.23 ng/ul	188392	Chrysene-d12	21.525

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	81
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81





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Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

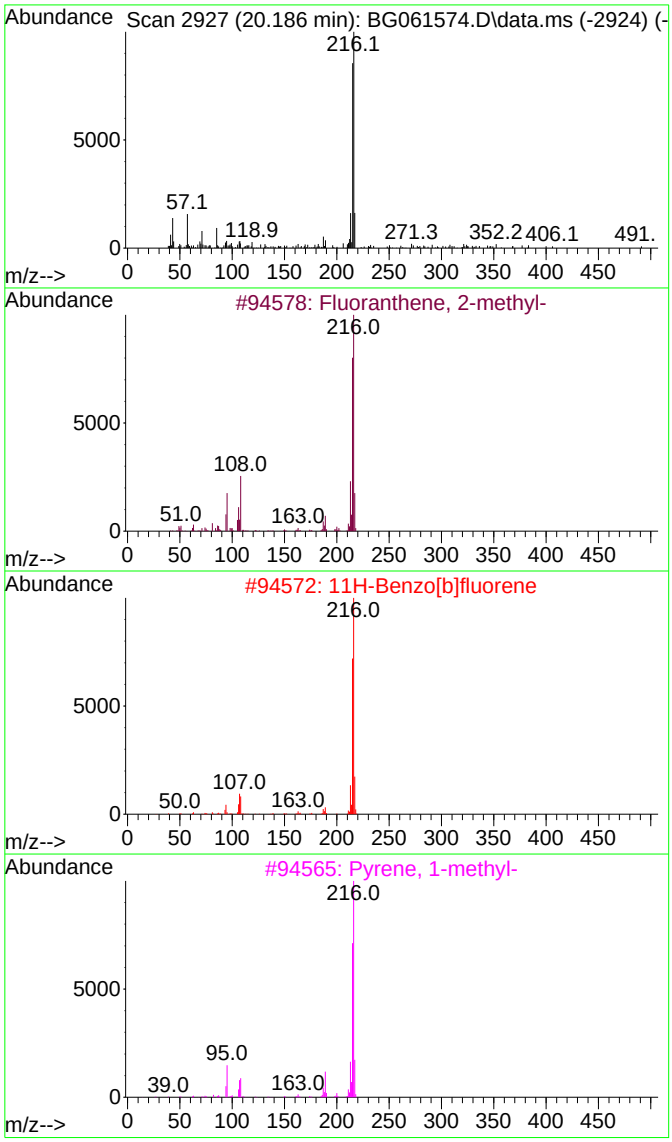
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

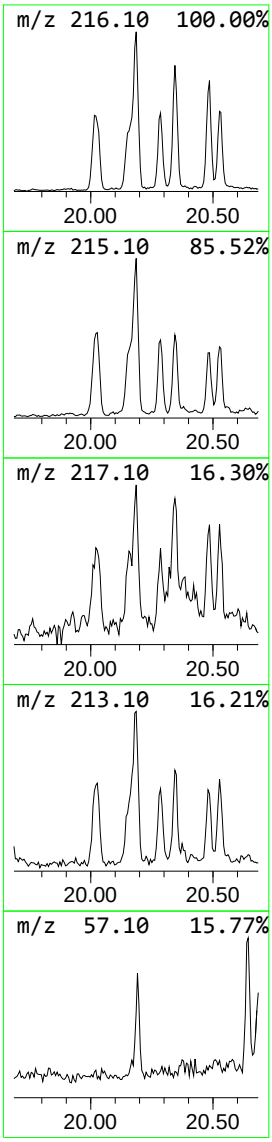
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Fluoranthene, 2-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.186	2.82 ng/ul	164886	Chrysene-d12	21.525

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	80
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	78
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	78
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5	1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	72





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Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

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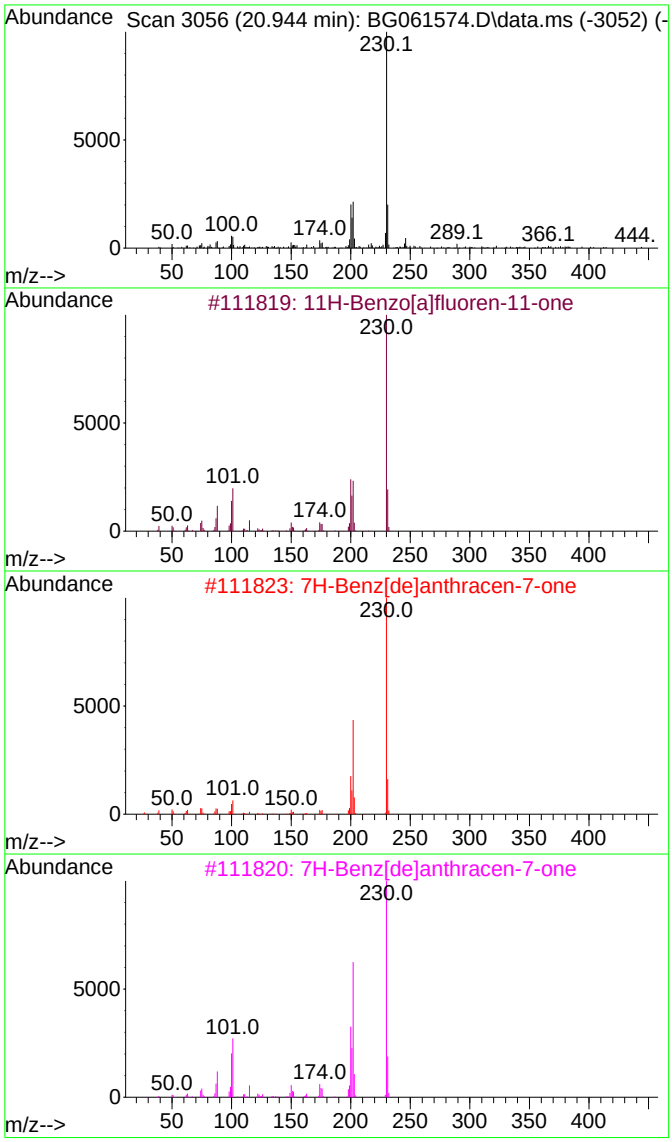
Quant Title : SVOA CALIBRATION

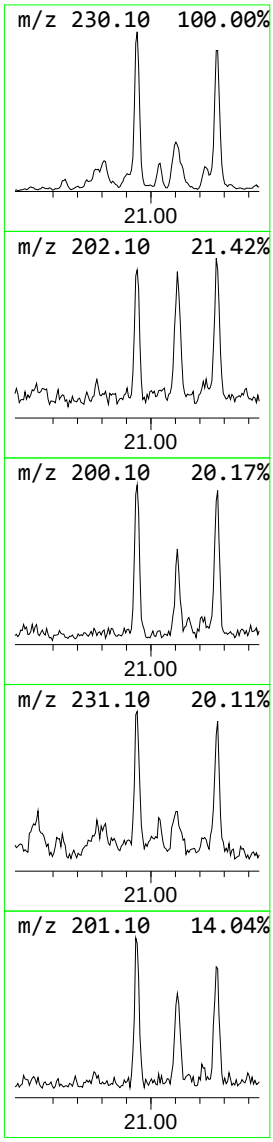
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 11H-Benzo[a]fluoren-11-one Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.944	2.48 ng/ul	144566	Chrysene-d12	21.525

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	96
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	89
3	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	76
5	1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	72





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Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

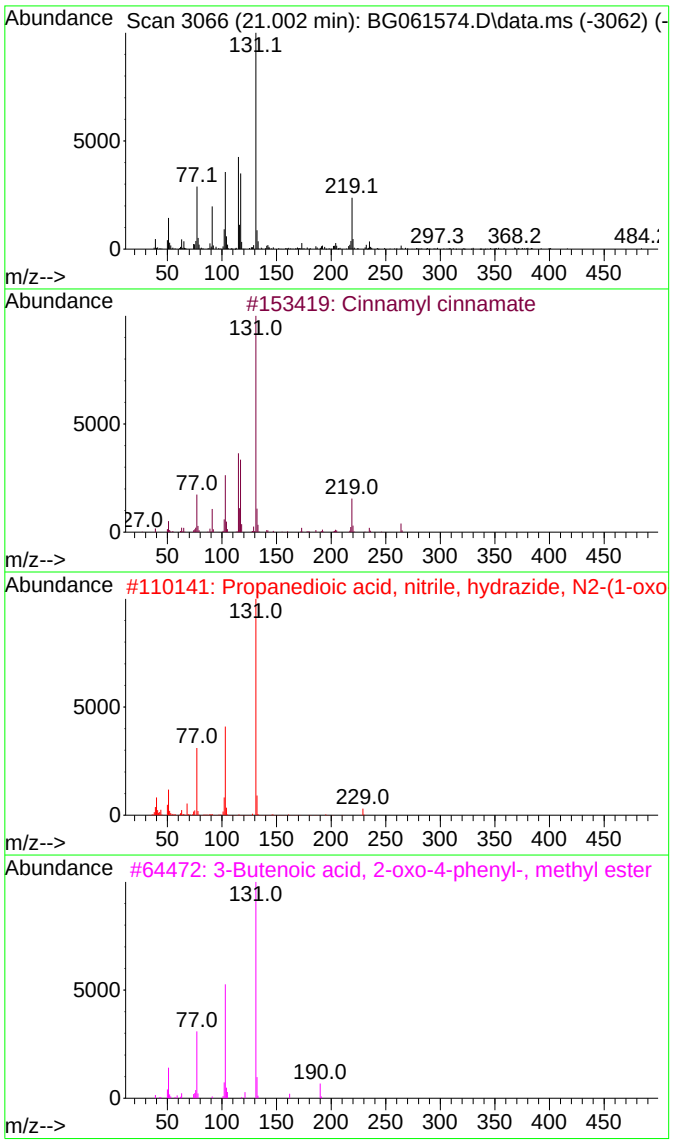
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

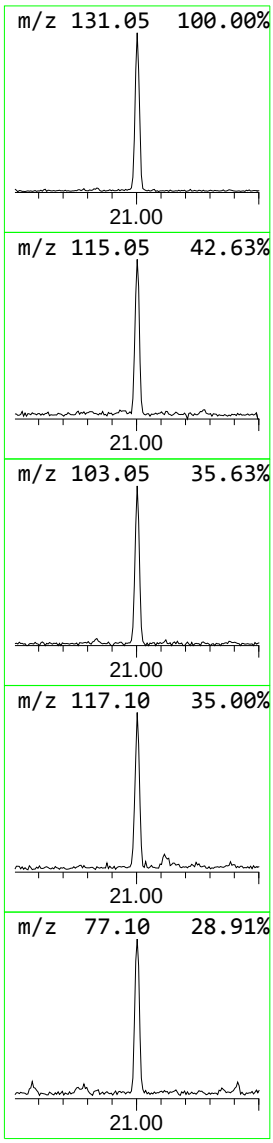
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Cinnamyl cinnamate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.002	4.51 ng/ul	263456	Chrysene-d12	21.525

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2	Propanedioic acid, nitrile, hydr...	229	C12H11N3O2	294878-35-6	46
3	3-Butenoic acid, 2-oxo-4-phenyl-...	190	C11H10O3	1000142-22-4	46
4	trans-1-Cinnamoylimidazole	198	C12H10N2O	001138-15-4	43
5	(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	43





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Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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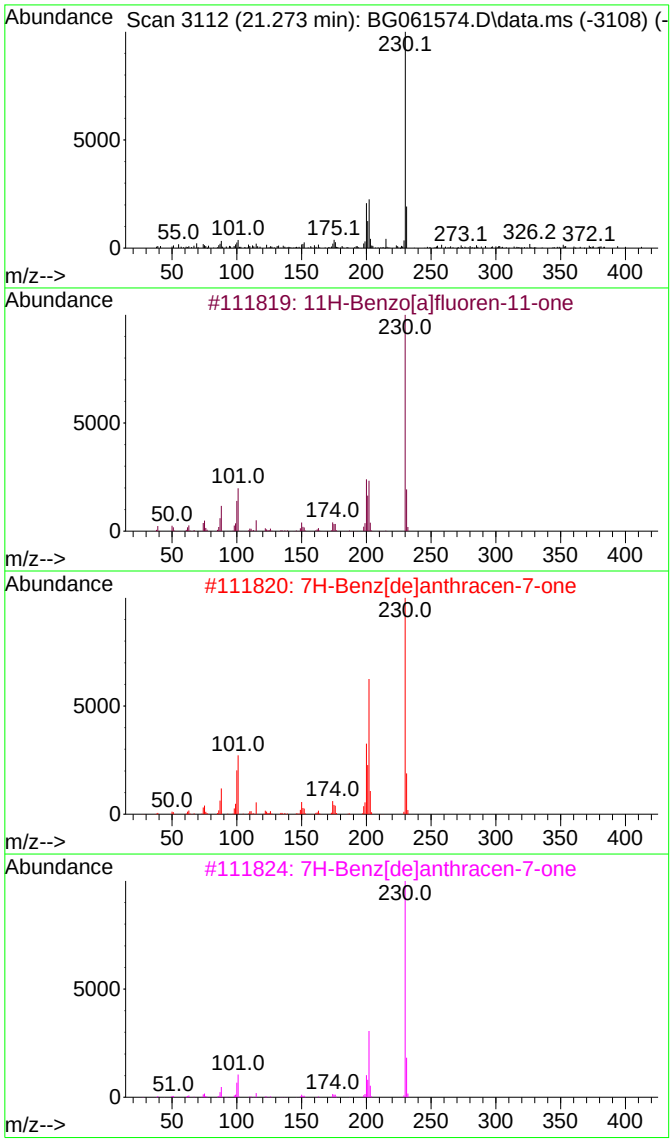
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

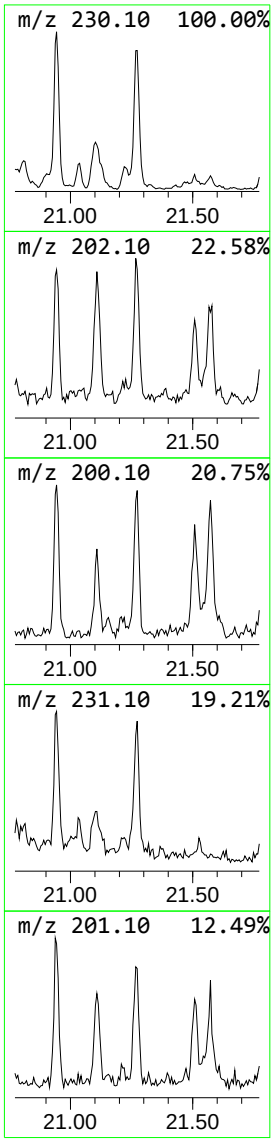
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 7H-Benz[de]anthracen-7-one Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.273	2.64 ng/ul	154305	Chrysene-d12	21.525

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	90
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
3	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	81
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	81
5	6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	81





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Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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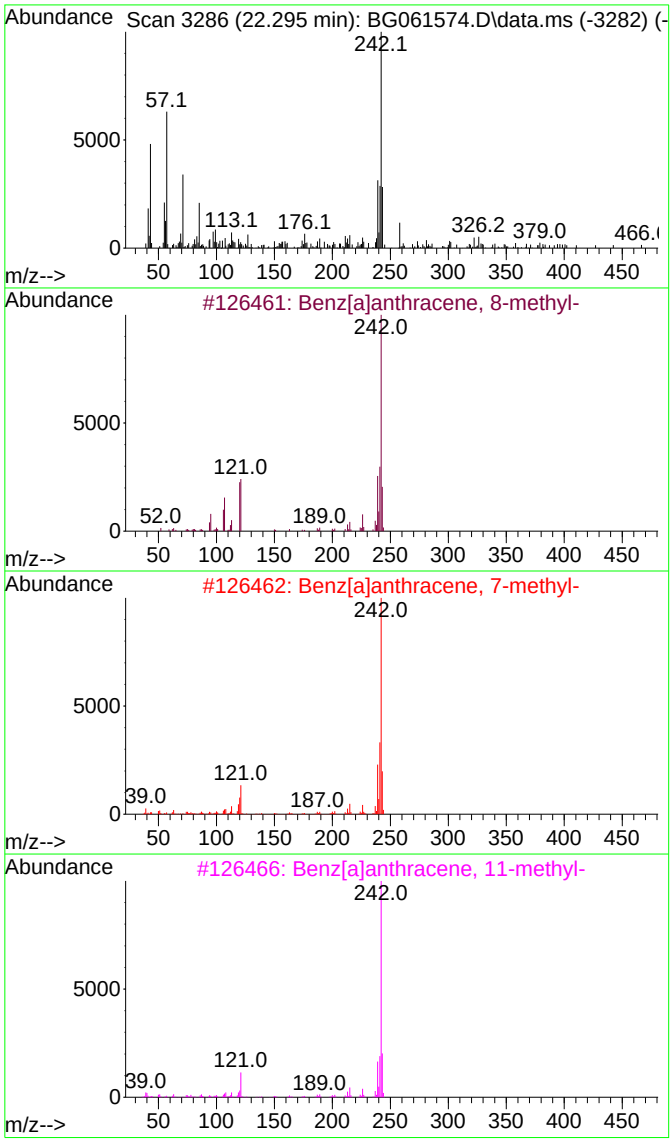
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

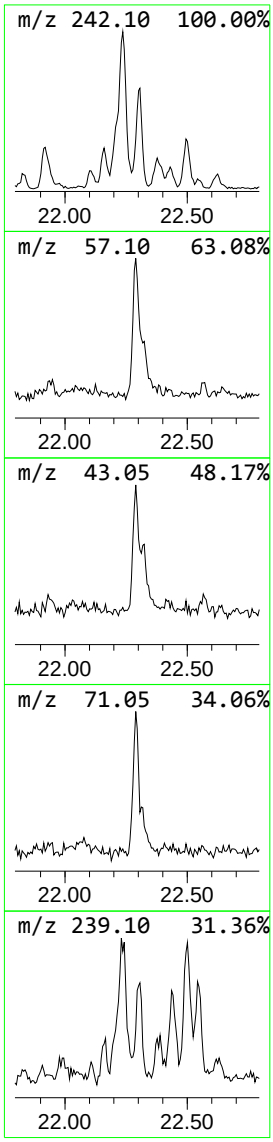
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Benz[a]anthracene, 8-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.295	3.97 ng/ul	231694	Chrysene-d12	21.525

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	64
2	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	60
3	Benz[a]anthracene, 11-methyl-	242	C19H14	006111-78-0	60
4	Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	60
5	Chrysene, 6-methyl-	242	C19H14	001705-85-7	58





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Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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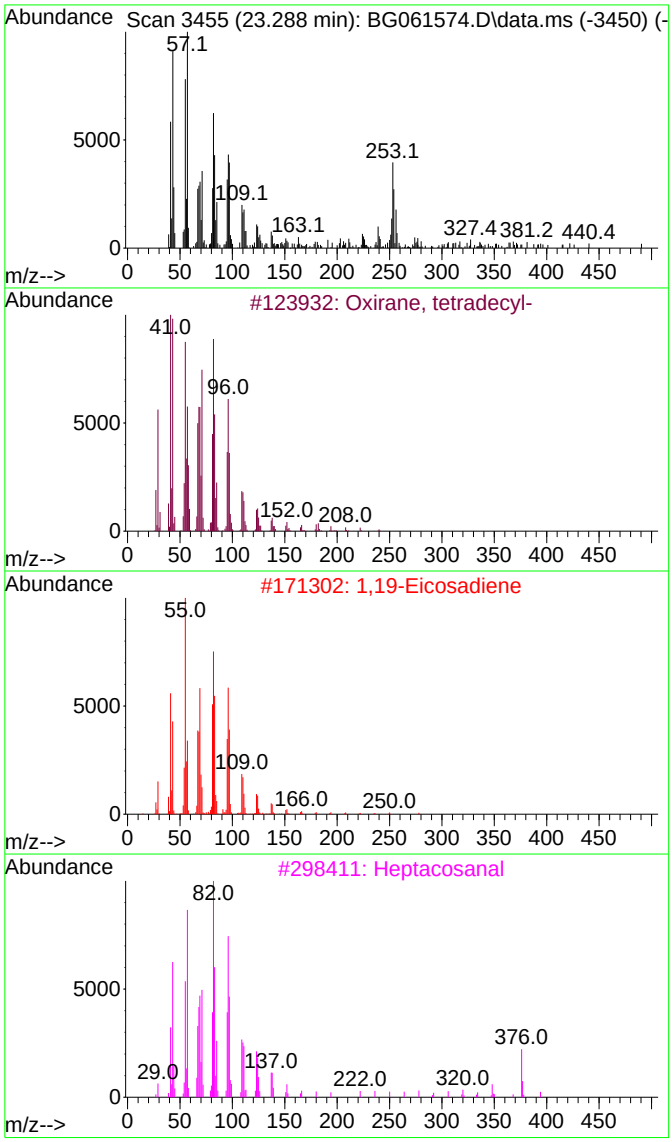
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

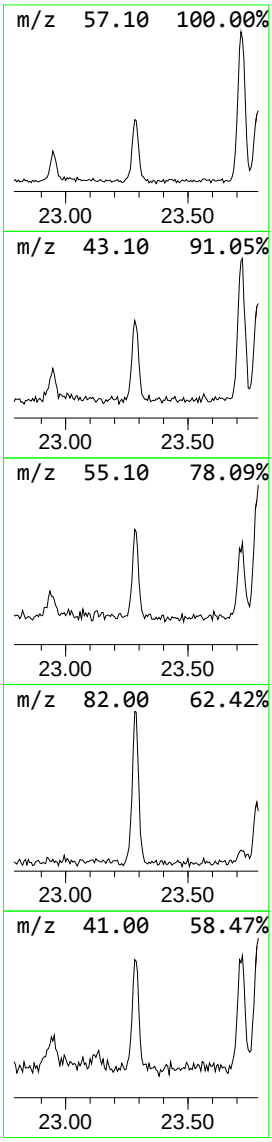
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Oxirane, tetradecyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.288	11.69 ng/ul	258598	Perylene-d12	24.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, tetradecyl-	240	C16H32O	007320-37-8	96
2	1,19-Eicosadiene	278	C20H38	014811-95-1	93
3	Heptacosanal	394	C27H54O	072934-03-3	92
4	Octadecanal	268	C18H36O	000638-66-4	89
5	Octadecanal	268	C18H36O	000638-66-4	78





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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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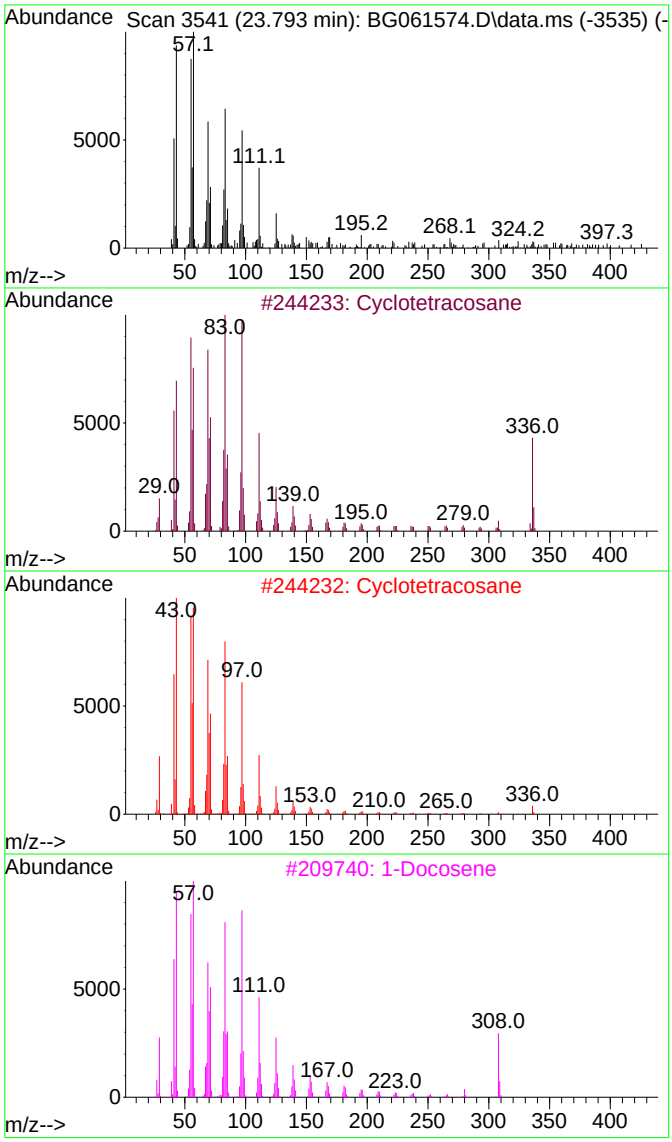
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

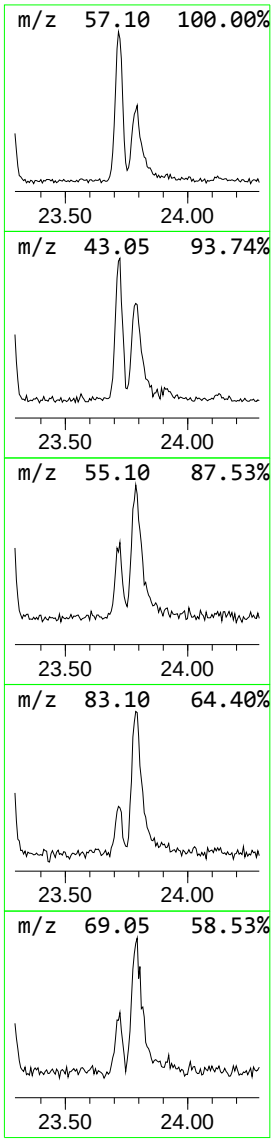
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Cyclic23.793 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.793	16.84 ng/ul	372415	Perylene-d12	24.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	95
2	Cyclotetracosane	336	C24H48	000297-03-0	95
3	1-Docosene	308	C22H44	001599-67-3	87
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	87
5	1-Methoxyhexacosane	396	C27H56O	237742-59-5	87





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Data File : BG061574.D
Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

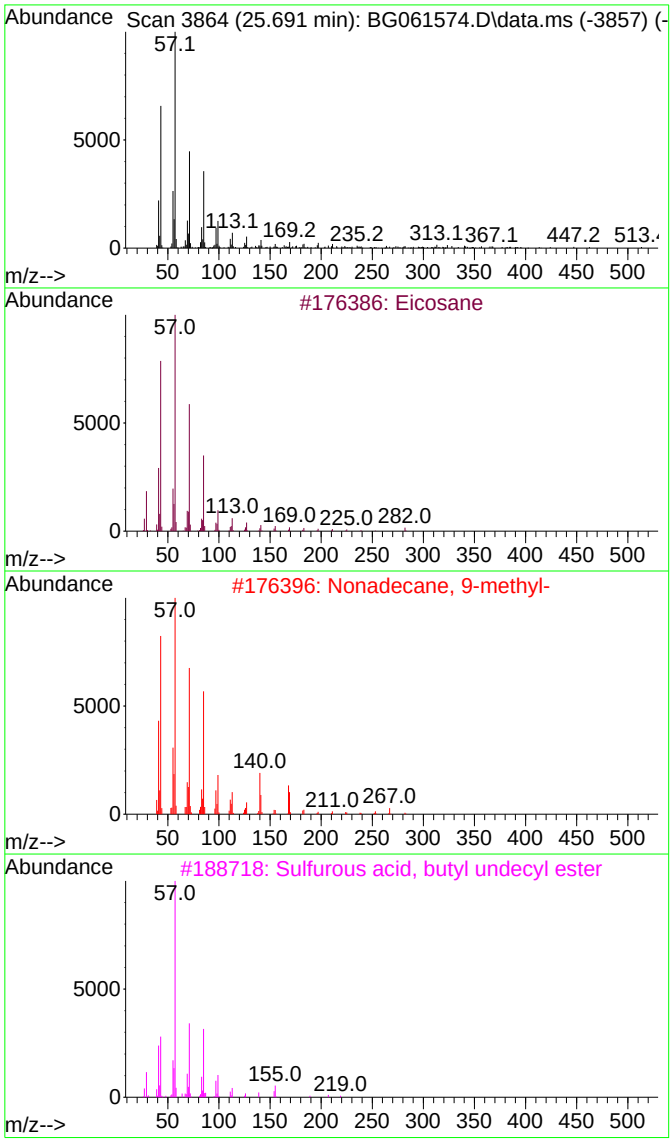
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

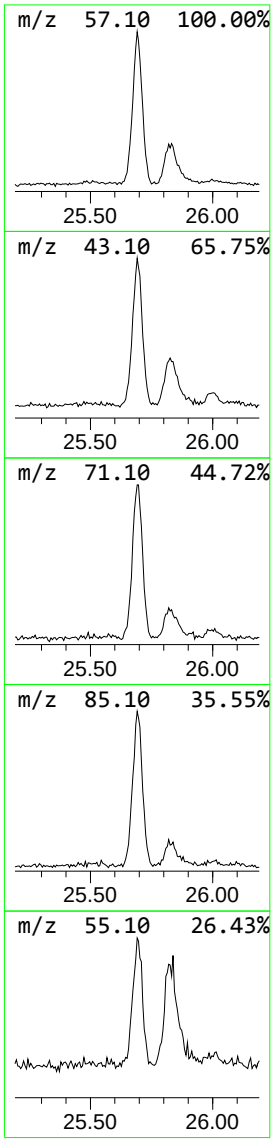
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.691	23.36 ng/ul	516656	Perylene-d12	24.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	90
2	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	87
3	Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	83
4	Heptadecane, 3-methyl-	254	C18H38	006418-44-6	83
5	Tetratetracontane	619	C44H90	007098-22-8	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
Data File : BG061574.D
Acq On : 8 Jun 2024 21:00
Operator : MA/JU
Sample : P2647-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BHBT1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.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...	3.076	177.5	ng/ul	1590370	1	7.871	179242	20.0
unknown-01	3.852	2.2	ng/ul	19609	1	7.871	179242	20.0
2-Pentanone, 4-...	4.992	3.0	ng/ul	26740	1	7.871	179242	20.0
Cyclopentene, 3...	6.555	3.3	ng/ul	29364	1	7.871	179242	20.0
Santolina triene	6.860	5.4	ng/ul	48031	1	7.871	179242	20.0
2,4,7,9-Tetrame...	13.411	24.0	ng/ul	407724	3	14.510	339822	20.0
Acetophenone, 4...	13.817	14.3	ng/ul	242170	3	14.510	339822	20.0
Dibenzofuran, 4...	15.855	2.2	ng/ul	36696	3	14.510	339822	20.0
(DEL) Alkane: S...	16.243	2.4	ng/ul	57372	4	17.265	475116	20.0
unknown-02	17.013	2.1	ng/ul	49250	4	17.265	475116	20.0
n-Hexadecanoic ...	18.165	13.2	ng/ul	312919	4	17.265	475116	20.0
Anthracene, 9-m...	18.194	7.2	ng/ul	169895	4	17.265	475116	20.0
Phenanthrene, 4...	18.276	2.1	ng/ul	50315	4	17.265	475116	20.0
unknown-03	18.335	8.5	ng/ul	201133	4	17.265	475116	20.0
1H-Cyclopropa[l...	18.376	4.0	ng/ul	95044	4	17.265	475116	20.0
5,16[1',2']:8,1...	18.640	3.6	ng/ul	86766	4	17.265	475116	20.0
9,10-Anthracene...	18.693	4.2	ng/ul	100056	4	17.265	475116	20.0
unknown-04	19.087	18.1	ng/ul	429298	4	17.265	475116	20.0
Cyclopenta(def)...	19.210	5.6	ng/ul	132784	4	17.265	475116	20.0
Pyrene, 1-methyl-	20.015	3.2	ng/ul	188392	5	21.525	1167900	20.0
Fluoranthene, 2...	20.186	2.8	ng/ul	164886	5	21.525	1167900	20.0
11H-Benzo[a]flu...	20.944	2.5	ng/ul	144566	5	21.525	1167900	20.0
Cinnamyl cinnamate	21.002	4.5	ng/ul	263456	5	21.525	1167900	20.0
7H-Benz[de]anth...	21.273	2.6	ng/ul	154305	5	21.525	1167900	20.0
Benz[a]anthrace...	22.295	4.0	ng/ul	231694	5	21.525	1167900	20.0
Oxirane, tetrad...	23.288	11.7	ng/ul	258598	6	24.633	442276	20.0
(DEL) Alkane: C...	23.793	16.8	ng/ul	372415	6	24.633	442276	20.0
(DEL) Alkane: S...	25.691	23.4	ng/ul	516656	6	24.633	442276	20.0