LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061348.D
Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BH658

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: BG061348.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.035	5	8	11	rVV2	12932	15204	1.31%	0.102%
2	3.082	11	16	31	rVB	267579	435005	37.46%	2.928%
3	3.335	56	59	64	rBV3	5295	6806	0.59%	0.046%
4	3.611	101	106	110	rBV	5487	7948	0.68%	0.053%
5	3.746	124	129	138	rBV2	22361	38689	3.33%	0.260%
6	4.304	221	224	229	rVB2	6171	7451	0.64%	0.050%
7	5.961	502	506	514	rVB2	4540	6827	0.59%	0.046%
8	7.066	682	694	705	rBV	121416	216679	18.66%	1.458%
9	7.213	712	719	728	rVB	78449	129842	11.18%	0.874%
10	7.418	747	754	761	rBV	119569	198447	17.09%	1.336%
11	7.882	825	833	839	rBV	122225	204153	17.58%	1.374%
12	7.941	839	843	849	rVB2	4979	6926	0.60%	0.047%
13	8.599	948	955	962	rBV	139443	235351	20.27%	1.584%
14	9.034	1022	1029	1036	rBV	127980	222411	19.15%	1.497%
15	9.592	1118	1124	1130	rVB2	17254	25752	2.22%	0.173%
16	9.757	1144	1152	1160	rBV	115279	200451	17.26%	1.349%
17	10.309	1239	1246	1256	rVB	161502	288344	24.83%	1.941%
18	10.673	1300	1308	1315	rBV	172639	308091	26.53%	2.073%
19	10.726	1315	1317	1321	rVB3	5063	6131	0.53%	0.041%
20	10.808	1325	1331	1340	rVV2	81769	155920	13.43%	1.049%
21	11.208	1393	1399	1403	rBV3	5046	6900	0.59%	0.046%
22	11.414	1430	1434	1440	rBV3	15790	26119	2.25%	0.176%
23	11.702	1478	1483	1487	rBV2	3171	4908	0.42%	0.033%
24	12.101	1544	1551	1556	rBV	6114	9036	0.78%	0.061%
25	12.330	1585	1590	1595	rVB4	5026	8322	0.72%	0.056%
26	12.553	1621	1628	1632	rBV4	5730	8354	0.72%	0.056%
27	13.341	1757	1762	1768	rBV2	11262	16547	1.43%	0.111%
28	13.687	1814	1821	1826	rVB4	3162	7380	0.64%	0.050%
29	13.834	1841	1846	1852	rBV2	6419	10450	0.90%	0.070%
30	13.922	1852	1861	1867	rVV	300510	429984	37.03%	2.894%
31	13.999	1871	1874	1879	rVB2	5018	6167	0.53%	0.042%
32	14.204	1900	1909	1923	rBV	317066	511264	44.03%	3.441%
33	14.416	1940	1945	1951	rVB3	10925	16441	1.42%	0.111%
34	14.516	1951	1962	1968	rBV	279422	443912	38.23%	2.988%
35	14.575	1968	1972	1982	rVB5	10484	20204	1.74%	0.136%
36	14.745	1991	2001	2017	rBV4	69513	167604	14.43%	1.128%

LSC Area Percent Report

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Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BH658

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	14.868	2020	2022	2026	rVV5	7153	7857	0.68%	0.053%
38	14.915	2026	2030	2035	rVB2	14450	20539	1.77%	0.138%
39	14.992	2038	2043	2047	rVB3	4325	7939	0.68%	0.053%
40	15.127	2064	2066	2072	rVB2	4044	6620	0.57%	0.045%
41	15.303	2089	2096	2100	rBV5	5811	10136	0.87%	0.068%
42	15.368	2103	2107	2112	rVB2	11189	12499	1.08%	0.084%
43	15.509	2124	2131	2136	rBV2	409667	613935	52.87%	4.132%
44	15.562	2136	2140	2144	rVB3	12410	17047	1.47%	0.115%
45	15.644	2148	2154	2160	rBV	99399	153740	13.24%	1.035%
46	15.773	2173	2176	2182	rVB5	6384	8438	0.73%	0.057%
47	15.855	2186	2190	2194	rBV3	6996	11566	1.00%	0.078%
48	15.996	2209	2214	2215	rBV3	6275	6048	0.52%	0.041%
49	16.085	2223	2229	2230	rBV4	3510	5162	0.44%	0.035%
50	16.232	2249	2254	2264	rBV5	19986	45021	3.88%	0.303%
51	16.572	2308	2312	2317	rVB5	11039	16653	1.43%	0.112%
52	16.613	2317	2319	2324	rVB4	5717	8565	0.74%	0.058%
53	16.801	2347	2351	2354	rBV3	9109	11820	1.02%	0.080%
54	16.837	2354	2357	2361	rVB4	4269	6836	0.59%	0.046%
55	16.919	2367	2371	2378	rVB4	18430	29058	2.50%	0.196%
56	16.995	2378	2384	2385	rBV4	7553	9981	0.86%	0.067%
57	17.019	2385	2388	2393	rVV3	16274	23539	2.03%	0.158%
58	17.078	2395	2398	2407	rVB3	15357	25313	2.18%	0.170%
59	17.266	2423	2430	2433	rBV	342631	525672	45.27%	3.538%
60	17.307	2433	2437	2441	rVV	239271	351981	30.31%	2.369%
61	17.360	2441	2446	2457	rVB	440513	681516	58.69%	4.587%
62	17.454	2458	2462	2468	rBV7	7019	12035	1.04%	0.081%
63	17.665	2495	2498	2504	rVB	19156	24166	2.08%	0.163%
64	17.759	2511	2514	2518	rBV5	11827	21656	1.87%	0.146%
65	17.841	2525	2528	2533	rVB6	12210	16019	1.38%	0.108%
66	18.059	2561	2565	2569	rBV5	9102	14494	1.25%	0.098%
67	18.106	2569	2573	2574	rBV3	11431	12111	1.04%	0.082%
68	18.141	2575	2579	2581	rBV	41955	61158	5.27%	0.412%
69	18.335	2606	2612	2615	rBV2	61689	115783	9.97%	0.779%
70	18.452	2628	2632	2636	rBV5	19313	29751	2.56%	0.200%
71	18.488	2636	2638	2643	rVV5	8808	14189	1.22%	0.095%
72	18.640	2660	2664	2667	rBV	40248	51587	4.44%	0.347%
73	18.682	2667	2671	2678	rVB2	50370	74558	6.42%	0.502%
74	18.858	2699	2701	2703	rBV3	7710	8218	0.71%	0.055%
75	18.887	2703	2706	2711	rVB4	19885	21458	1.85%	0.144%
76	18.934	2712	2714	2718	rVB2	16752	18900	1.63%	0.127%

LSC Area Percent Report

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Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BH658

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

77	18.975	2718	2721	2726	rVB4	13937	17457	1.50%	0.117%
78	19.058	2730	2735	2738	rBV	47167	78303	6.74%	0.527%
79	19.199	2755	2759	2768	rVB3	38321	77866	6.71%	0.524%
80	19.322	2774	2780	2785	rVB	597131	813487	70.06%	5.475%
81	19.416	2793	2796	2799	rBV3	14773	23840	2.05%	0.160%
82	19.657	2831	2837	2839	rBV2	603638	945092	81.39%	6.361%
83	19.686	2839	2842	2849	rVB	523758	711929	61.31%	4.791%
84	19.757	2850	2854	2858	rBV3	35762	49666	4.28%	0.334%
85	20.004	2891	2896	2903	rBV3	47695	123609	10.65%	0.832%
86	20.274	2939	2942	2946	rBV	36917	51173	4.41%	0.344%
87	20.474	2973	2976	2979	rBV	41605	51224	4.41%	0.345%
88	20.521	2981	2984	2988	rVB	42933	55732	4.80%	0.375%
89	20.685	3010	3012	3015	rVB	38104	36722	3.16%	0.247%
90	20.932	3050	3054	3061	rVB	80772	122123	10.52%	0.822%
91	21.185	3093	3097	3104	rVB3	97971	183926	15.84%	1.238%
92	21.408	3132	3135	3139	rVB	117439	147291	12.68%	0.991%
93	21.514	3145	3153	3157	rBV2	508493	1161167	100.00%	7.815%
94	21.561	3158	3161	3171	rVB	301213	453795	39.08%	3.054%
95	21.707	3183	3186	3190	rVB2	53438	67864	5.84%	0.457%
96	22.289	3282	3285	3291	rVB4	63863	100244	8.63%	0.675%
97	23.282	3450	3454	3465	rVB3	110233	215008	18.52%	1.447%
98	23.623	3506	3512	3521	rBV	216602	728046	62.70%	4.900%
99	24.393	3637	3643	3649	rBV3	246837	559303	48.17%	3.764%
100	24.604	3672	3679	3686	rBV2	223524	560051	48.23%	3.769%

Sum of corrected areas: 14858502

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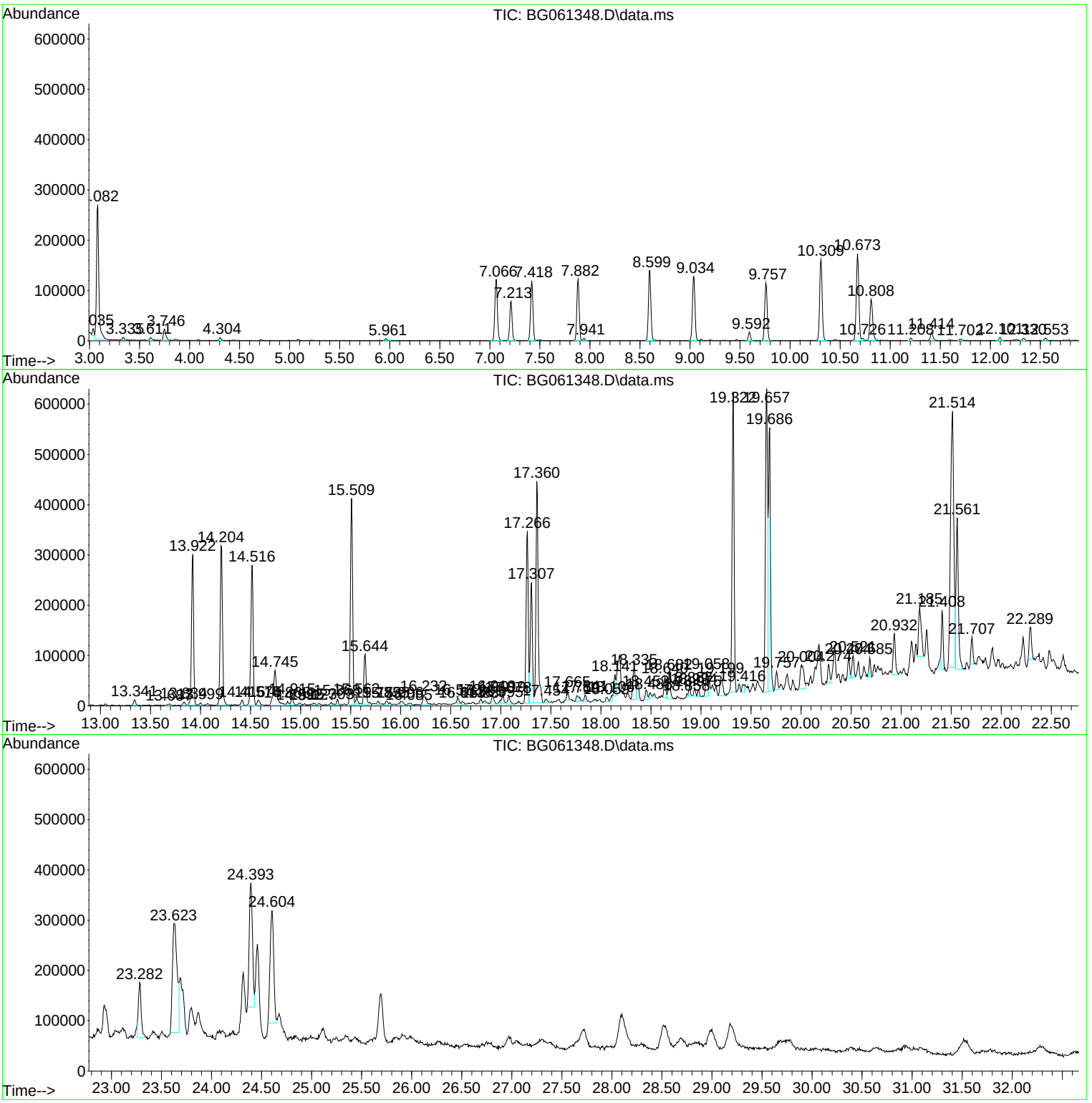
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

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\BG052924\
Data File : BG061348.D
Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

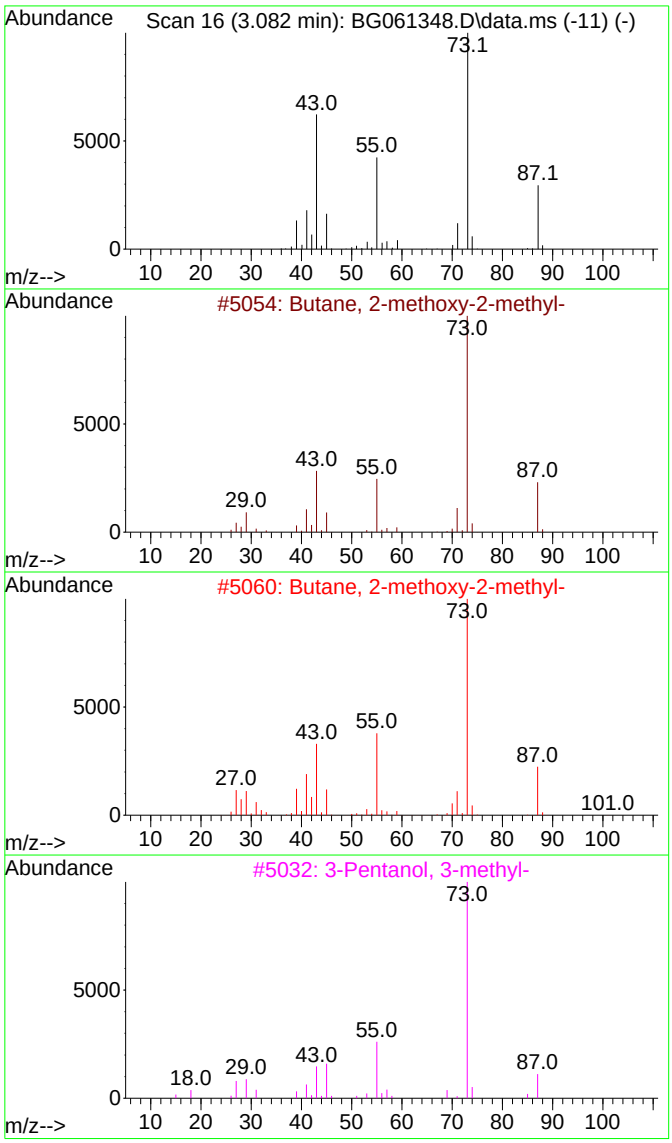
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

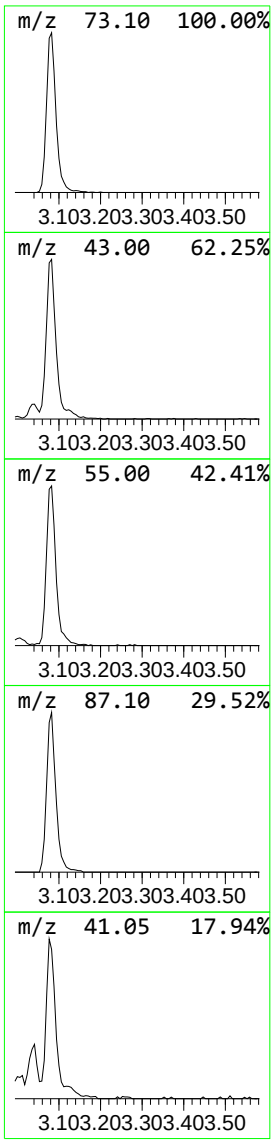
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.082	42.62 ng/ul	435005	1,4-Dichlorobenzene-d4	7.877

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28
5	4-Penten-2-ol, acetate	128	C7H12O2	1000352-31-5	10





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061348.D
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Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

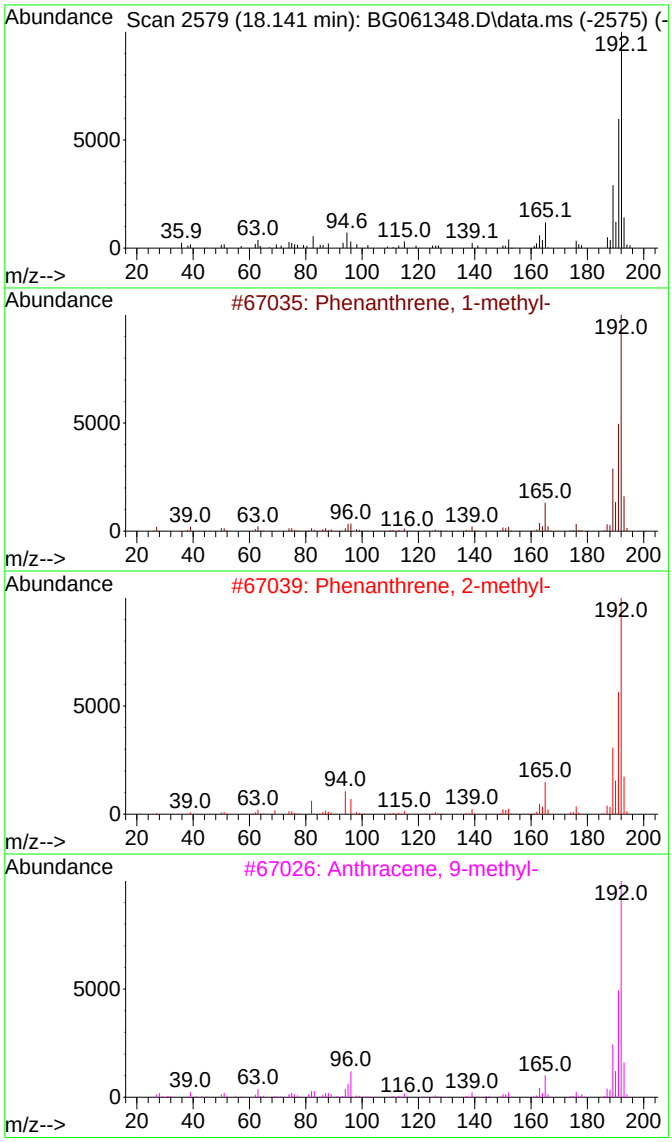
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

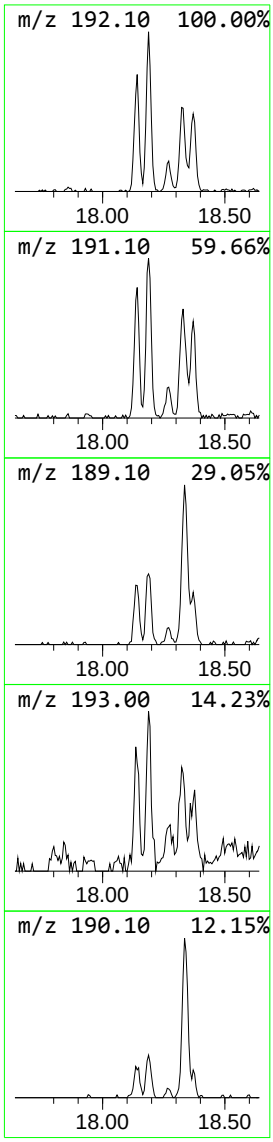
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.141	2.33 ng/ul	61158	Phenanthrene-d10	17.266

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

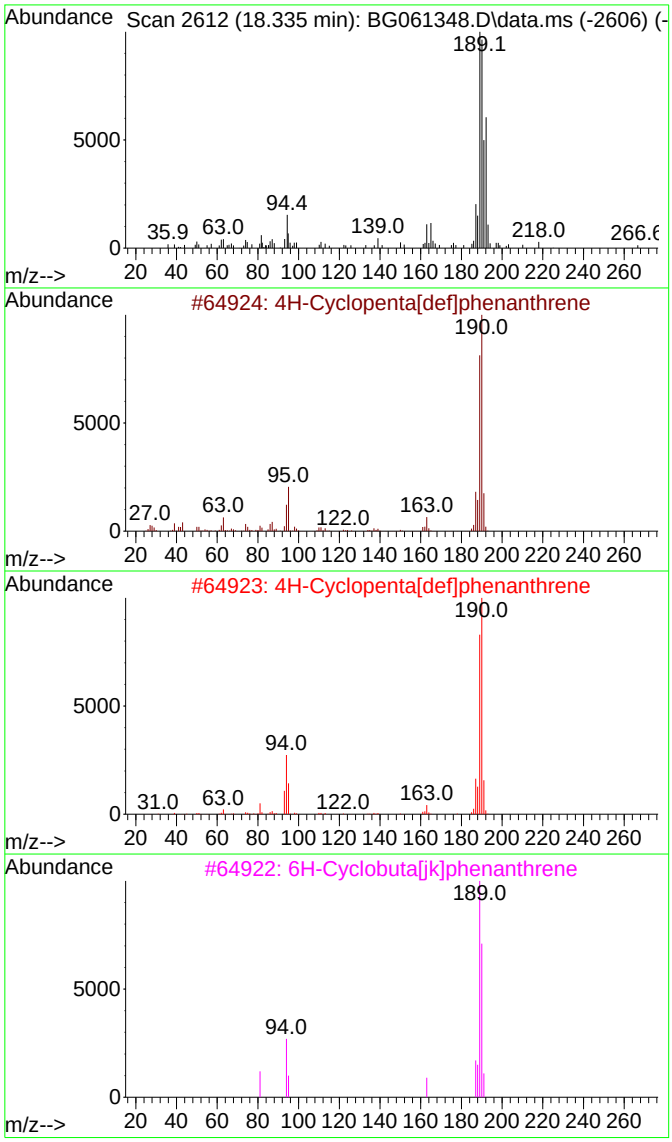
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

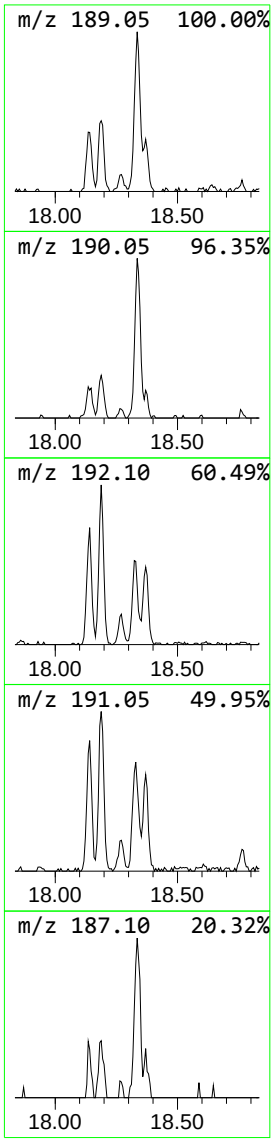
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.335	4.41 ng/ul	115783	Phenanthrene-d10	17.266

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
5	Isoquinoline, 1-chloro-3-ethyl-	191	C11H10ClN	055150-52-2	38





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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

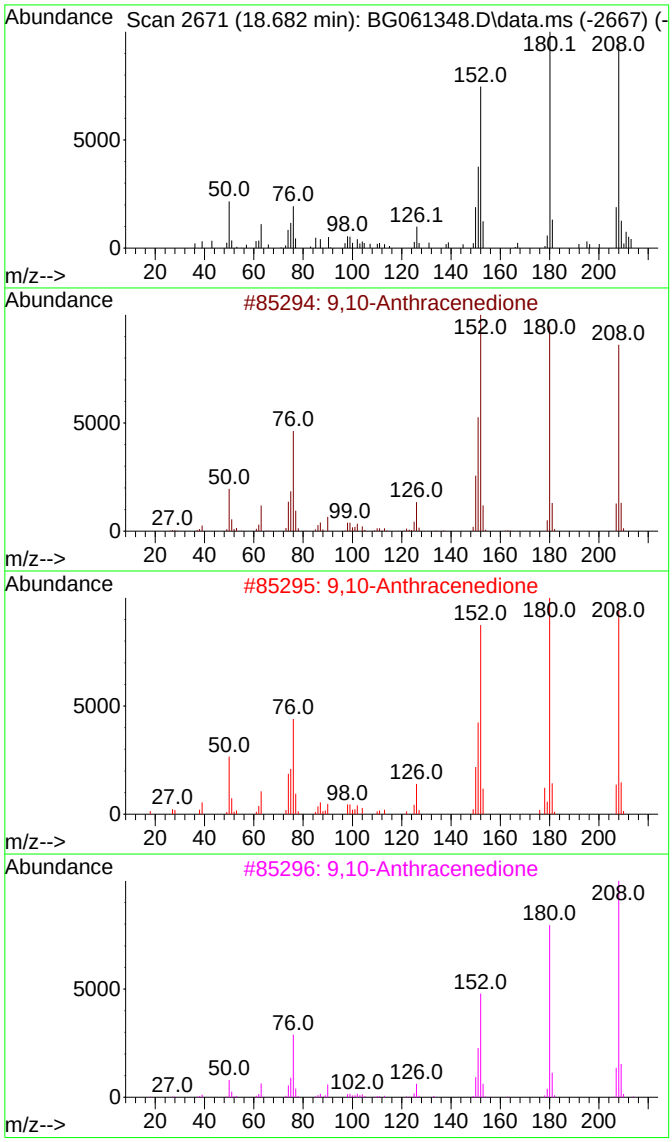
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

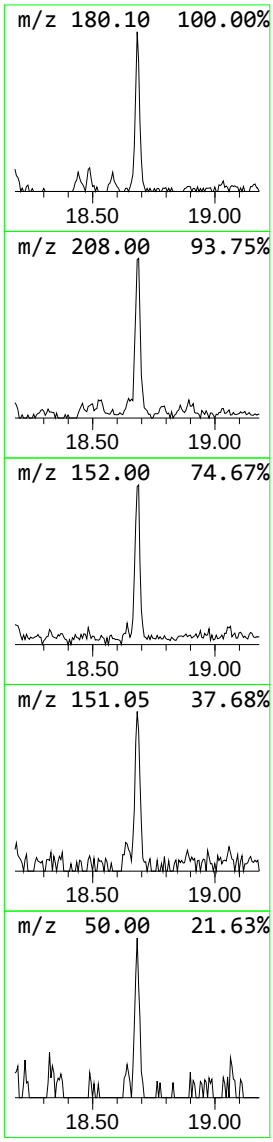
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.682	2.84 ng/ul	74558	Phenanthrene-d10	17.266

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061348.D
Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

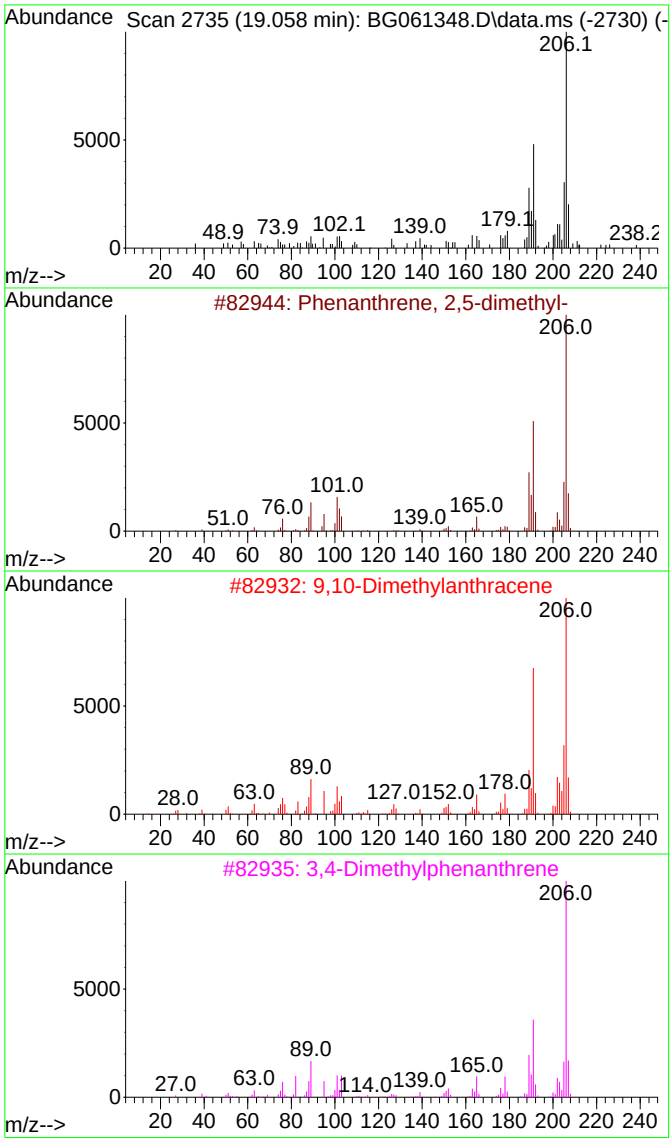
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

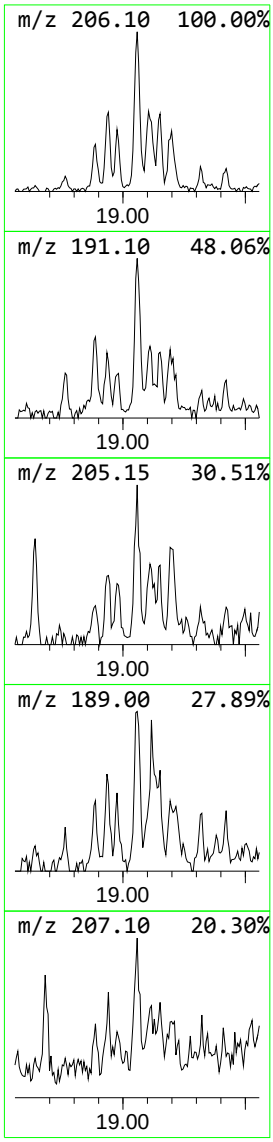
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.058	2.98 ng/ul	78303	Phenanthrene-d10	17.266

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
3	3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	90
4	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	87
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	87





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Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

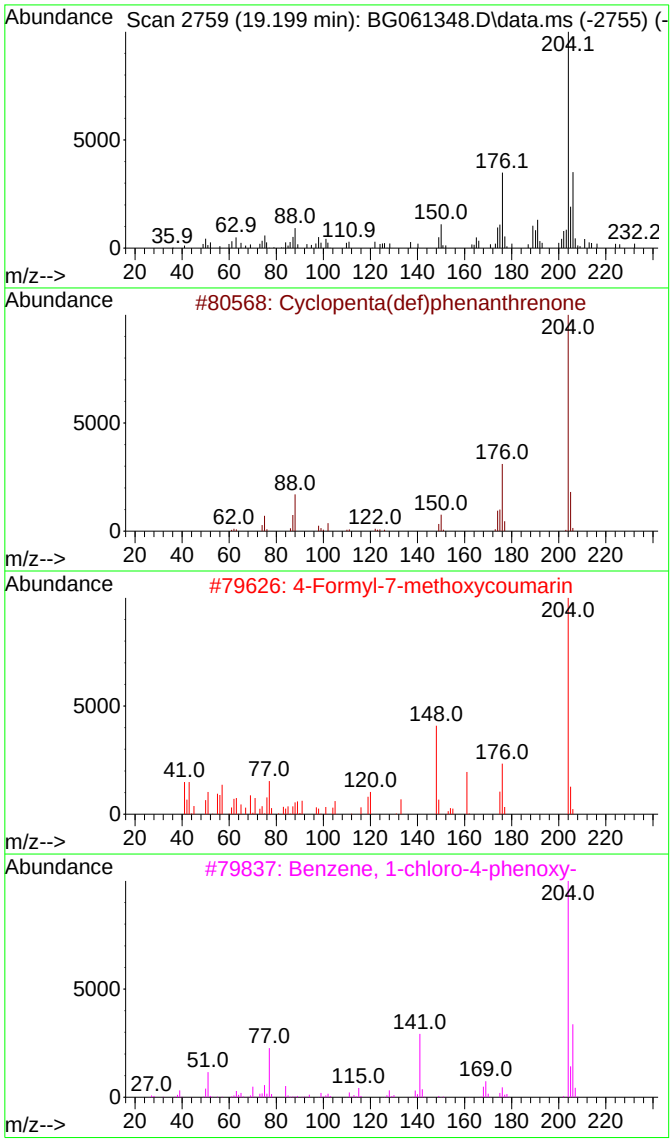
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

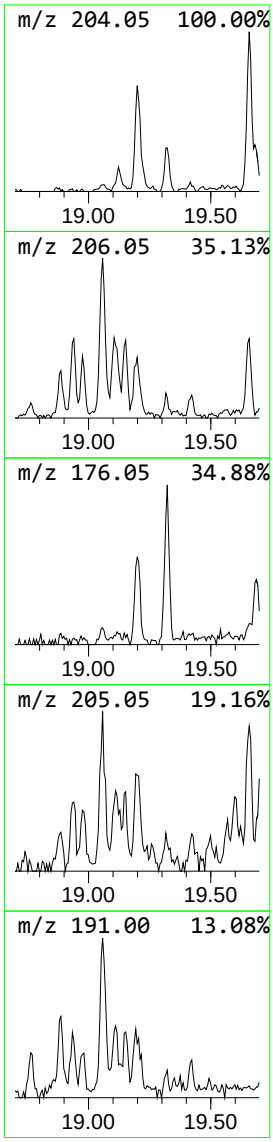
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclopenta(def)phenanthrenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.199	2.96 ng/ul	77866	Phenanthrene-d10	17.266

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	76
2	4-Formyl-7-methoxycoumarin	204	C11H8O4	1000429-03-0	50
3	Benzene, 1-chloro-4-phenoxy-	204	C12H9ClO	007005-72-3	49
4	[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	49
5	[1,1'-Biphenyl-4-ol], 2-chloro-	204	C12H9ClO	023719-22-4	47





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Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

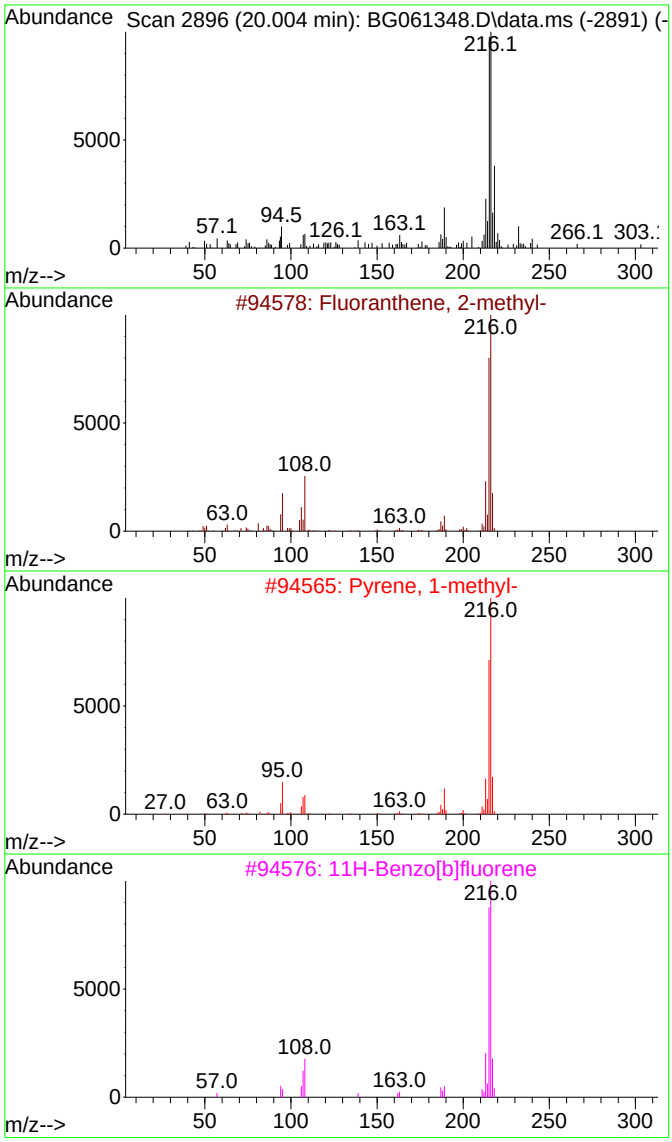
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

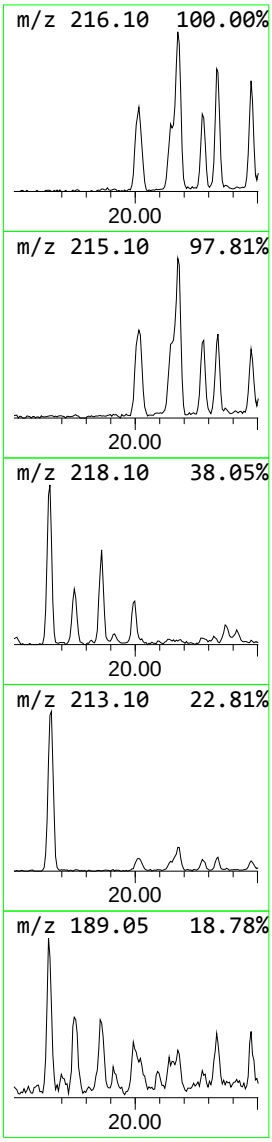
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.004	2.13 ng/ul	123609	Chrysene-d12	21.514

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	70
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	60
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	58
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	58
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	55





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Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

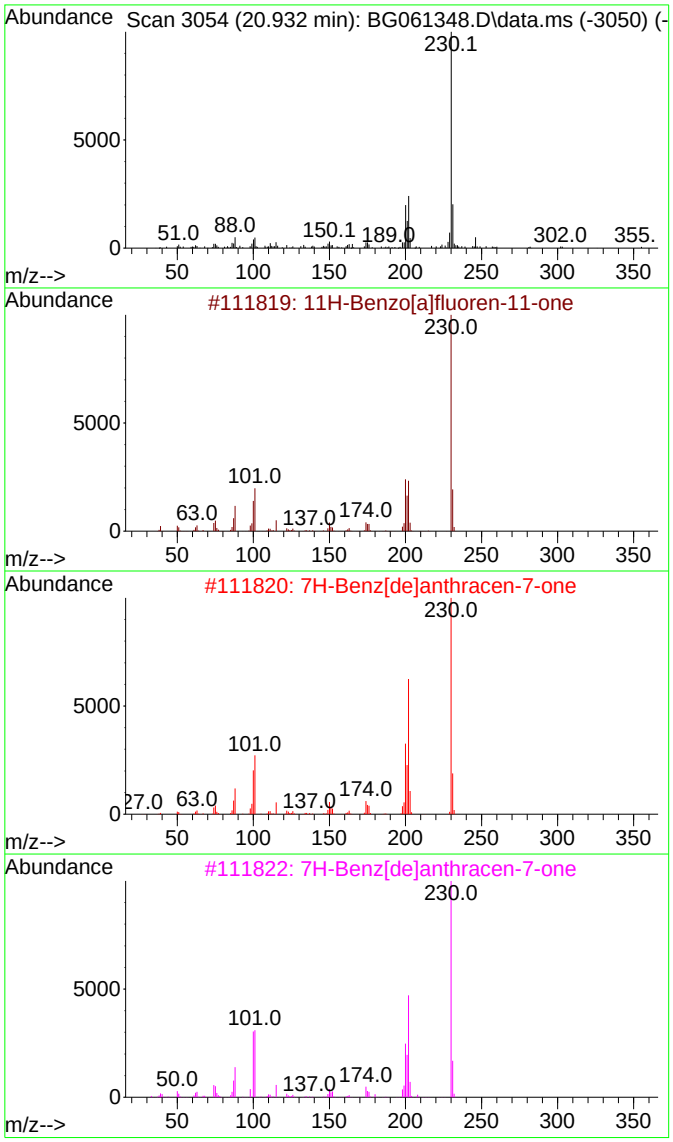
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

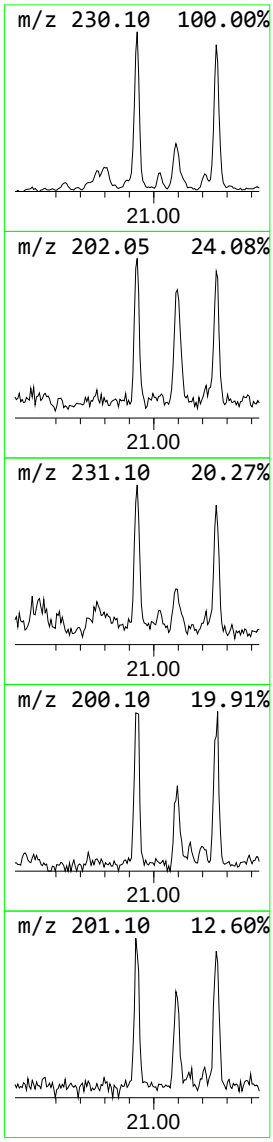
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 11H-Benzo[a]fluoren-11-one Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.932	2.10 ng/ul	122123	Chrysene-d12	21.514

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	93
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
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	81
5	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	74





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Data File : BG061348.D
Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

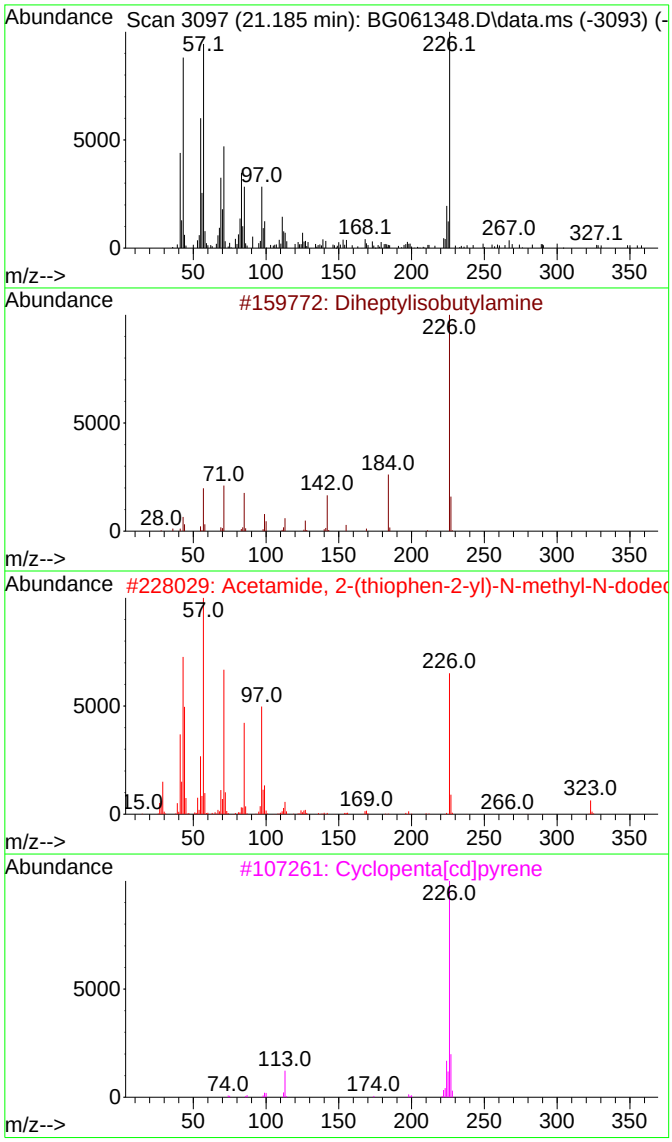
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

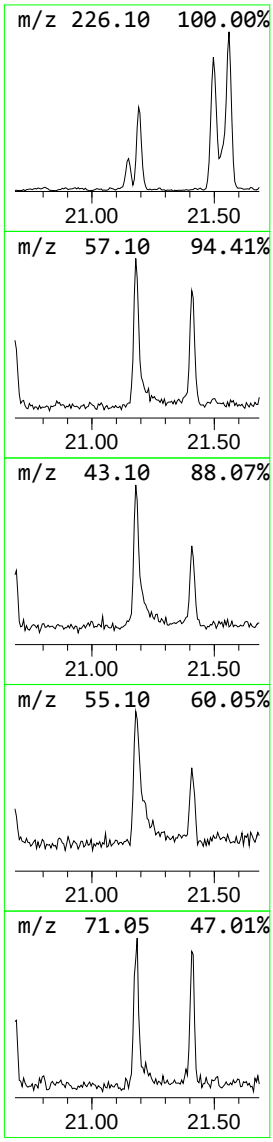
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Diheptylisobutylamine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.185	3.17 ng/ul	183926	Chrysene-d12	21.514

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diheptylisobutylamine	269	C18H39N	1000310-32-0	50
2	Acetamide, 2-(thiophen-2-yl)-N-m...	323	C19H33NOS	1000421-20-9	47
3	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	32
4	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	27
5	Heptadecane, 3-methyl-	254	C18H38	006418-44-6	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061348.D
Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

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

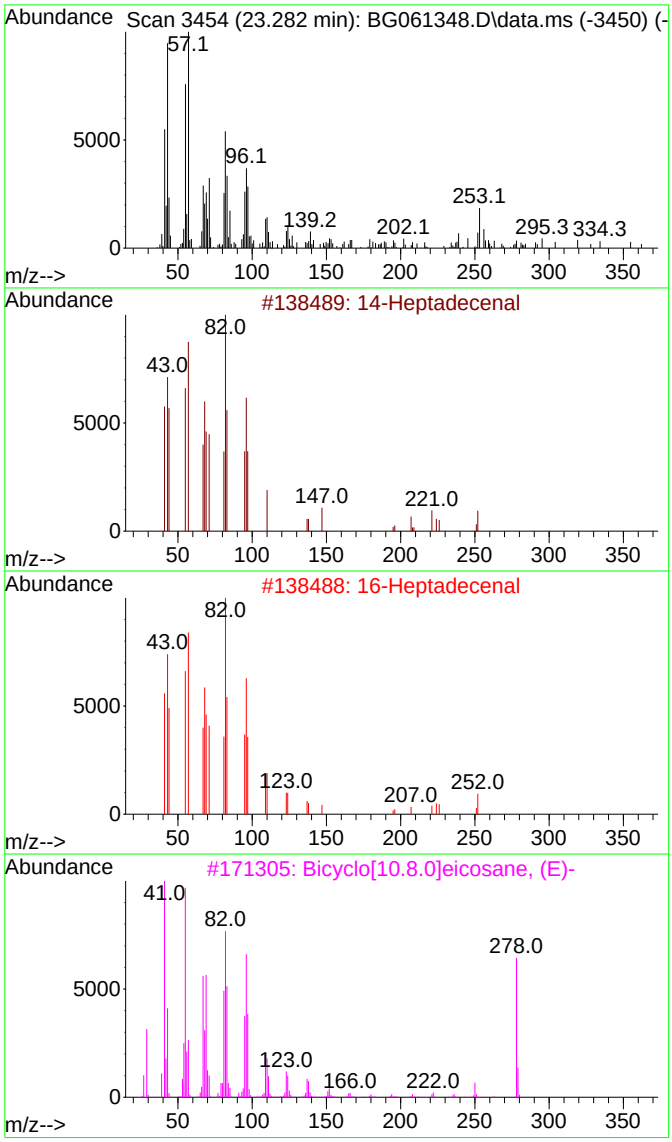
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

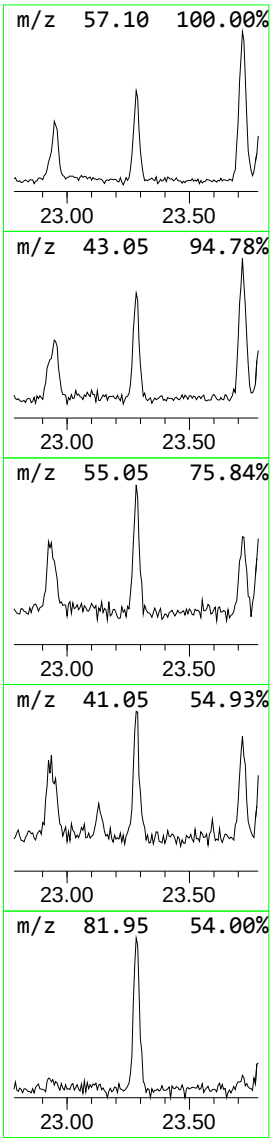
****

Peak Number 10 14-Heptadecenal Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.282	7.68 ng/ul	215008	Perylene-d12	24.598

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	14-Heptadecenal	252	C17H32O	1010144-58-0	90
2	16-Heptadecenal	252	C17H32O	1010144-57-9	90
3	Bicyclo[10.8.0]eicosane, (E)-	278	C20H38	1010155-85-0	90
4	1,19-Eicosadiene	278	C20H38	014811-95-1	90
5	Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	90





Tentatively Identified Compound (LSC) summary

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Data File : BG061348.D
Acq On : 30 May 2024 4:59
Operator : MA/JU
Sample : P2571-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH658

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...	3.082	42.6	ng/ul	435005	1	7.877	204153	20.0
Phenanthrene, 1...	18.141	2.3	ng/ul	61158	4	17.266	525672	20.0
4H-Cyclopenta[d...	18.335	4.4	ng/ul	115783	4	17.266	525672	20.0
9,10-Anthracene...	18.682	2.8	ng/ul	74558	4	17.266	525672	20.0
Phenanthrene, 2...	19.058	3.0	ng/ul	78303	4	17.266	525672	20.0
Cyclopenta(def)...	19.199	3.0	ng/ul	77866	4	17.266	525672	20.0
Fluoranthene, 2...	20.004	2.1	ng/ul	123609	5	21.514	1161170	20.0
11H-Benzo[a]flu...	20.932	2.1	ng/ul	122123	5	21.514	1161170	20.0
Diheptylisobuty...	21.185	3.2	ng/ul	183926	5	21.514	1161170	20.0
14-Heptadecenal	23.282	7.7	ng/ul	215008	6	24.598	560051	20.0