LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A4D39

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_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP021323.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	27	30	38	rVB2	15865	25752	0.42%	0.031%
2	3.675	113	117	123	rVB	42028	58735	0.96%	0.071%
3	4.781	299	305	315	rBV	26723	47653	0.78%	0.058%
4	6.858	651	658	673	rBV	129098	374860	6.14%	0.455%
5	6.981	673	679	696	rVB	142269	300819	4.92%	0.365%
6	7.181	706	713	725	rBV	204233	405766	6.64%	0.492%
7	7.622	781	788	801	rBV	502444	865264	14.17%	1.049%
8	8.375	909	916	928	rBV	161765	440876	7.22%	0.535%
9	8.793	979	987	1007	rBV	170711	369366	6.05%	0.448%
10	9.369	1079	1085	1094	rBV	6577	21512	0.35%	0.026%
11	9.505	1101	1108	1130	rBV	132091	310588	5.08%	0.377%
12	10.081	1198	1206	1230	rBV2	119861	471636	7.72%	0.572%
13	10.381	1250	1257	1274	rBV	591404	1216649	19.92%	1.475%
14	10.604	1286	1295	1308	rBV4	24398	92391	1.51%	0.112%
15	11.175	1384	1392	1393	rBV2	19184	30622	0.50%	0.037%
16	12.075	1539	1545	1549	rBV2	10502	18459	0.30%	0.022%
17	12.275	1576	1579	1584	rBV4	9175	15717	0.26%	0.019%
18	13.304	1749	1754	1762	rVB5	14905	31674	0.52%	0.038%
19	13.498	1781	1787	1797	rBV2	51970	179823	2.94%	0.218%
20	13.681	1812	1818	1829	rVV	455407	661065	10.82%	0.802%
21	13.951	1856	1864	1878	rBV2	511557	987624	16.17%	1.198%
22	14.075	1882	1885	1891	rVV8	12827	21656	0.35%	0.026%
23	14.263	1910	1917	1924	rBV	1069021	1630207	26.69%	1.977%
24	14.328	1924	1928	1935	rVB	64399	98206	1.61%	0.119%
25	14.487	1944	1955	1963	rBV5	115686	357824	5.86%	0.434%
26	14.628	1974	1979	1984	rBV2	81002	178713	2.93%	0.217%
27	14.681	1986	1988	1991	rVB	31041	34767	0.57%	0.042%
28	14.745	1997	1999	2003	rVB5	17864	18032	0.30%	0.022%
29	14.845	2012	2016	2021	rBV3	14287	23517	0.39%	0.029%
30	15.281	2084	2090	2097	rBV	626365	1108460	18.15%	1.344%
31	15.345	2097	2101	2109	rVB	62116	104439	1.71%	0.127%
32	15.475	2117	2123	2131	rBV3	108979	273657	4.48%	0.332%
33	15.545	2131	2135	2140	rVB4	35202	54836	0.90%	0.066%
34	15.640	2148	2151	2156	rVB	20333	30895	0.51%	0.037%
35	15.798	2174	2178	2184	rVB	20084	31551	0.52%	0.038%
36	16.028	2211	2217	2220	rBV8	12344	25715	0.42%	0.031%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

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_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

37	16.175	2238	2242	2246	rBV6	24188	34939	0.57%	0.042%
38	16.287	2257	2261	2267	rVB2	27342	38511	0.63%	0.047%
39	16.375	2270	2276	2281	rBV6	16415	43193	0.71%	0.052%
40	16.492	2290	2296	2311	rBV	860612	1709920	27.99%	2.073%
41	16.751	2337	2340	2345	rVB	44234	59837	0.98%	0.073%
42	16.810	2345	2350	2359	rBV5	85463	205786	3.37%	0.250%
43	16.898	2362	2365	2370	rVB	75214	96602	1.58%	0.117%
44	16.951	2371	2374	2375	rBV3	22265	25188	0.41%	0.031%
45	16.986	2376	2380	2387	rVB5	65101	113989	1.87%	0.138%
46	17.087	2388	2397	2401	rBV	1369048	2196684	35.96%	2.664%
47	17.128	2401	2404	2410	rVV	1141141	1721573	28.19%	2.088%
48	17.192	2410	2415	2419	rVV	845724	1309157	21.43%	1.587%
49	17.228	2419	2421	2425	rVV	224044	241942	3.96%	0.293%
50	17.263	2425	2427	2431	rVB4	36566	44921	0.74%	0.054%
51	17.539	2463	2474	2479	rBV2	156900	408046	6.68%	0.495%
52	17.622	2485	2488	2490	rBV2	26268	32658	0.53%	0.040%
53	17.698	2495	2501	2510	rVB	455875	795309	13.02%	0.964%
54	17.798	2516	2518	2528	rVB7	46501	103462	1.69%	0.125%
55	17.892	2529	2534	2539	rBV4	69195	142290	2.33%	0.173%
56	17.957	2540	2545	2548	rBV2	308586	453191	7.42%	0.550%
57	18.016	2549	2555	2558	rVV2	1164131	1717296	28.12%	2.082%
58	18.045	2558	2560	2564	rVB	227932	251399	4.12%	0.305%
59	18.122	2571	2573	2578	rVB	79112	105101	1.72%	0.127%
60	18.186	2579	2584	2589	rBV2	485396	831069	13.61%	1.008%
61	18.251	2592	2595	2598	rVB	237349	273900	4.48%	0.332%
62	18.298	2598	2603	2606	rBV2	192933	283338	4.64%	0.344%
63	18.386	2613	2618	2627	rBV	563295	1084242	17.75%	1.315%
64	18.475	2629	2633	2636	rBV3	146869	243418	3.99%	0.295%
65	18.586	2647	2652	2657	rBV2	203995	396445	6.49%	0.481%
66	18.669	2662	2666	2670	rVB	137361	196726	3.22%	0.239%
67	18.816	2689	2691	2695	rBV	42457	43859	0.72%	0.053%
68	18.986	2717	2720	2724	rVB2	191261	254011	4.16%	0.308%
69	19.045	2725	2730	2733	rBV	326580	482693	7.90%	0.585%
70	19.081	2733	2736	2740	rVB	319319	386171	6.32%	0.468%
71	19.222	2752	2760	2768	rBV	3383849	5099467	83.49%	6.184%
72	19.345	2778	2781	2787	rVB4	398672	602606	9.87%	0.731%
73	19.575	2815	2820	2822	rBV2	1716632	2296971	37.61%	2.785%
74	19.604	2822	2825	2829	rVB	2278029	2846181	46.60%	3.451%
75	19.675	2835	2837	2842	rBV	84387	100539	1.65%	0.122%
76	19.828	2861	2863	2867	rVB	126960	143152	2.34%	0.174%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

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_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

77	20.122	2910	2913	2917	rVB3	466439	660049	10.81%	0.800%
78	20.228	2927	2931	2935	rBV	240912	352770	5.78%	0.428%
79	20.369	2953	2955	2960	rVB3	131110	150497	2.46%	0.182%
80	20.480	2971	2974	2980	rVB	336457	456033	7.47%	0.553%
81	20.704	3008	3012	3017	rBV	339375	586283	9.60%	0.711%
82	21.104	3077	3080	3083	rBV	333615	417546	6.84%	0.506%
83	21.139	3084	3086	3088	rVV	345157	414606	6.79%	0.503%
84	21.175	3088	3092	3104	rVB5	1038707	2336152	38.25%	2.833%
85	21.422	3132	3134	3142	rVB	385747	572562	9.37%	0.694%
86	21.533	3145	3153	3158	rVV2	2261293	6108109	100.00%	7.407%
87	21.575	3158	3160	3168	rVB	1506773	2130472	34.88%	2.583%
88	21.727	3184	3186	3191	rBV2	237819	270483	4.43%	0.328%
89	21.916	3209	3218	3228	rVB3	1298216	3804181	62.28%	4.613%
90	22.357	3288	3293	3297	rBV	709448	1267364	20.75%	1.537%
91	22.398	3297	3300	3313	rVB3	1027351	2202616	36.06%	2.671%
92	23.421	3470	3474	3480	rVB3	606153	1230863	20.15%	1.493%
93	23.786	3529	3536	3543	rBV	1183738	3363523	55.07%	4.079%
94	23.980	3562	3569	3578	rBV2	1140270	3080783	50.44%	3.736%
95	24.598	3670	3674	3681	rBV	442163	1098361	17.98%	1.332%
96	24.674	3683	3687	3697	rVB	740575	1855632	30.38%	2.250%
97	24.827	3706	3713	3724	rVB	1094494	3062401	50.14%	3.713%
98	25.386	3802	3808	3818	rVB	953813	2262066	37.03%	2.743%
99	26.004	3905	3913	3924	rVB	1434368	4404691	72.11%	5.341%
100	26.180	3937	3943	3958	rVB2	591148	2045730	33.49%	2.481%

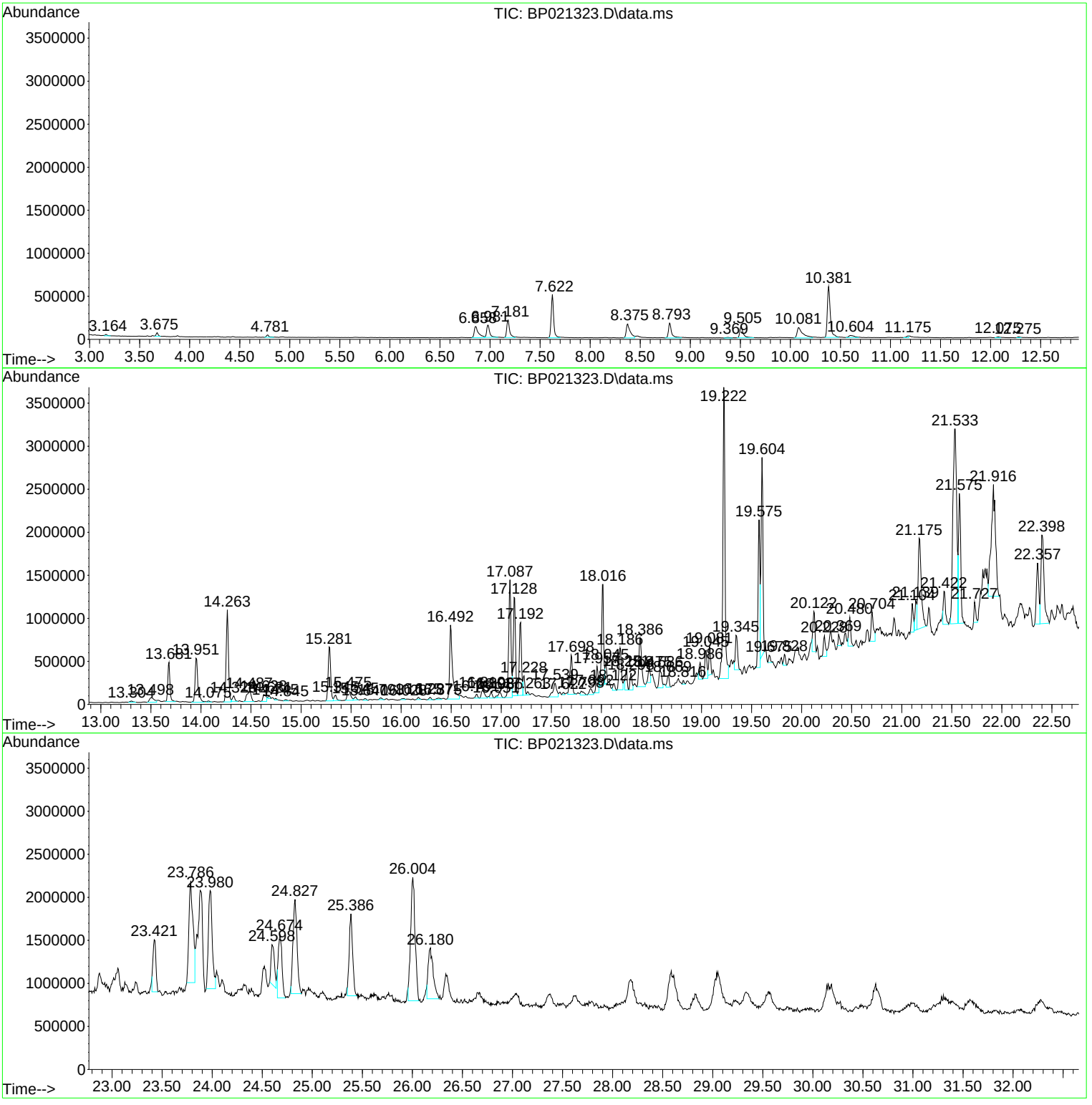
Sum of corrected areas: 82468881

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

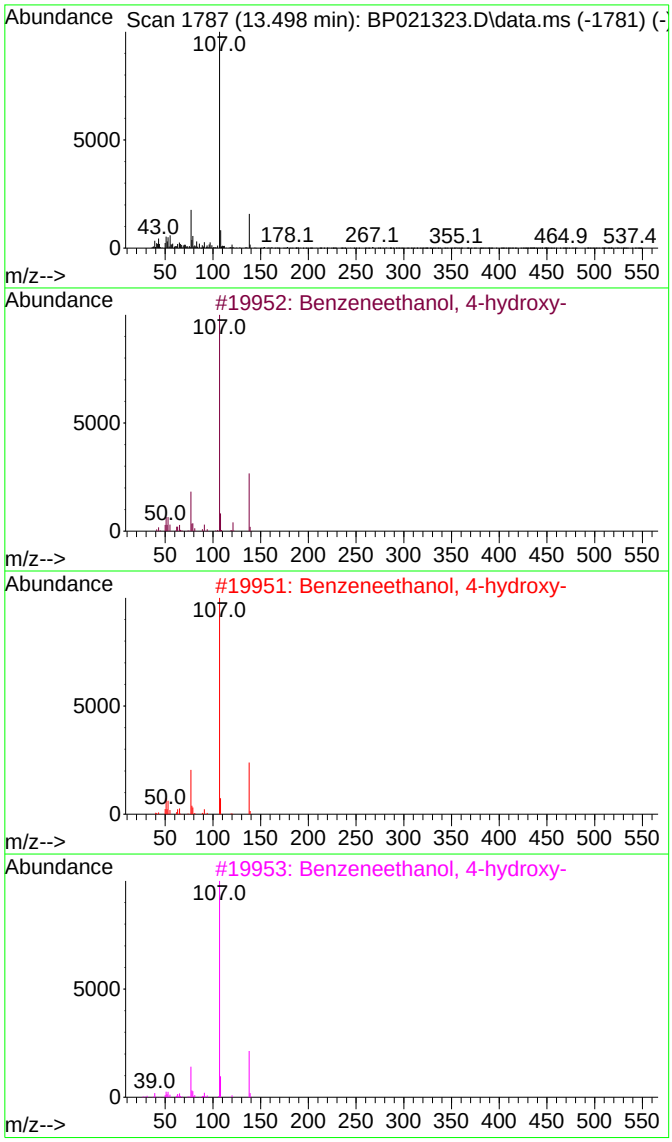
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

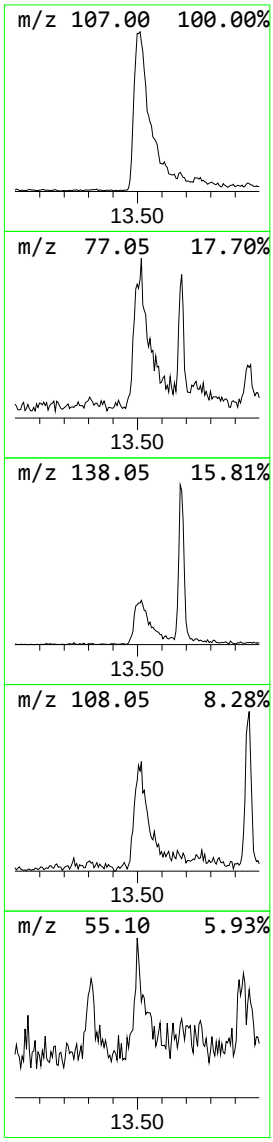
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzeneethanol, 4-hydroxy- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.499	2.21 ng/ul	179823	Acenaphthene-d10	14.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneethanol, 4-hydroxy-	138	C8H10O2	000501-94-0	87
2	Benzeneethanol, 4-hydroxy-	138	C8H10O2	000501-94-0	87
3	Benzeneethanol, 4-hydroxy-	138	C8H10O2	000501-94-0	87
4	Benzeneethanol, 4-hydroxy-	138	C8H10O2	000501-94-0	86
5	Benzeneethanol, 4-(acetyloxy)-	180	C10H12O3	060037-43-6	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

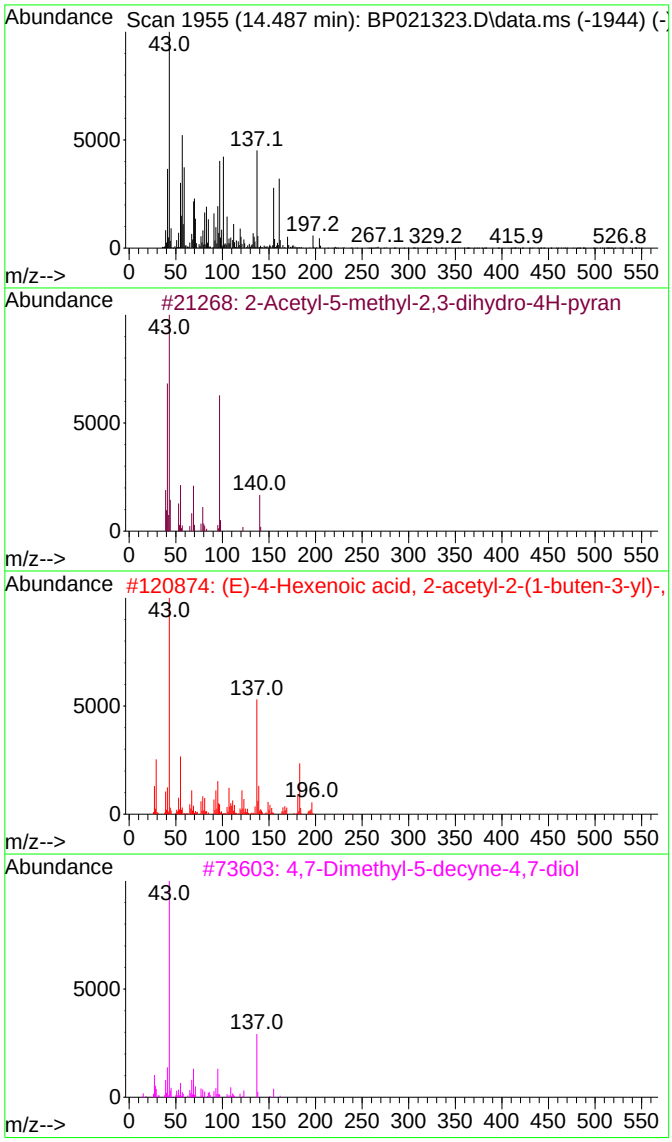
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

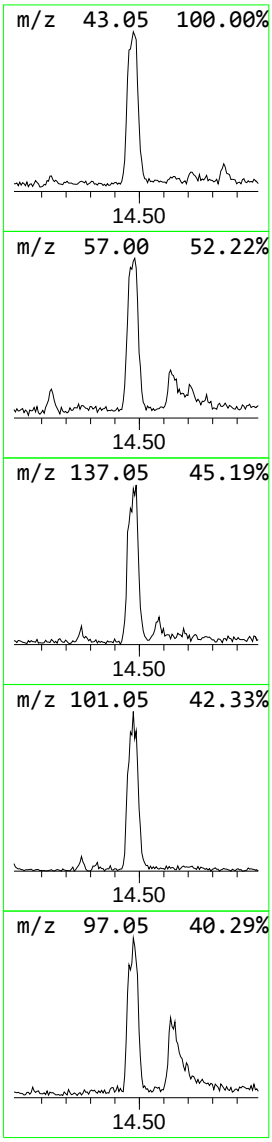
****

Peak Number 2 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.487	4.39 ng/ul	357824	Acenaphthene-d10	14.263

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Acetyl-5-methyl-2,3-dihydro-4H...	140	C8H12O2	1000288-90-9	10
2	(E)-4-Hexenoic acid, 2-acetyl-2-...	238	C14H22O3	1000159-09-7	10
3	4,7-Dimethyl-5-decyne-4,7-diol	198	C12H22O2	000126-87-4	10
4	4-Hepten-3-one, 5-methyl-, (E)-	126	C8H14O	020685-44-3	9
5	Pentanoic acid, 4,4-dimethyl-3-o...	158	C8H14O3	055107-14-7	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

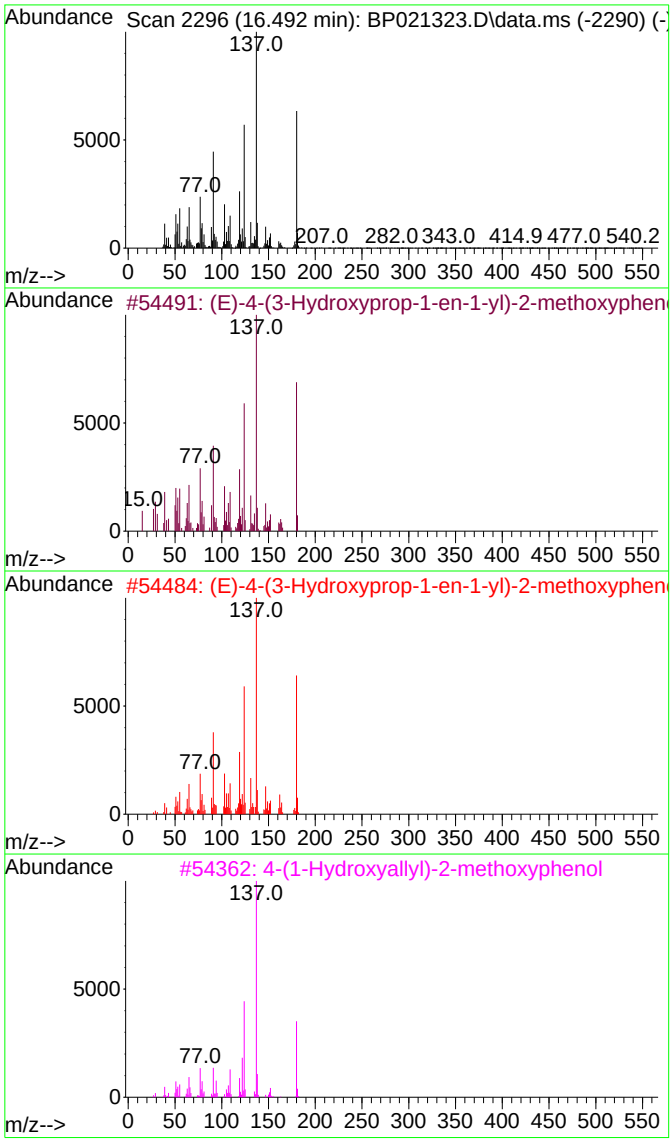
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

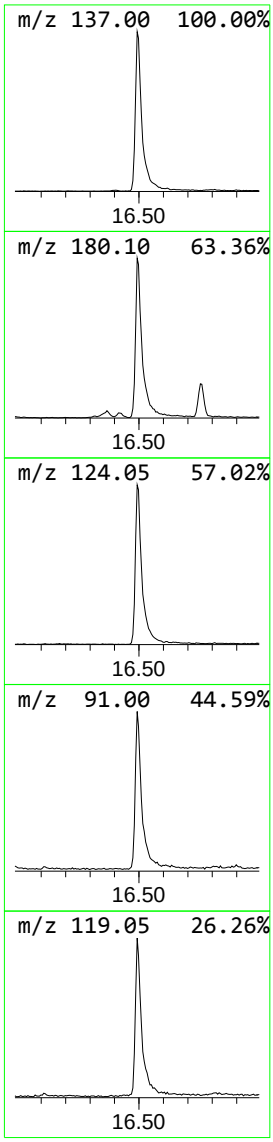
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (E)-4-(3-Hydroxyprop-1-en-1... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.492	15.57 ng/ul	1709920	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-4-(3-Hydroxyprop-1-en-1-yl)-...	180	C10H12O3	032811-40-8	99
2	(E)-4-(3-Hydroxyprop-1-en-1-yl)-...	180	C10H12O3	032811-40-8	95
3	4-(1-Hydroxyallyl)-2-methoxyphenol	180	C10H12O3	112465-50-6	68
4	1-Hydroxy-1-(4-methoxyphenyl)pro...	180	C10H12O3	015482-29-8	49
5	Guaiacol, 4-butyl-	180	C11H16O2	059832-96-1	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

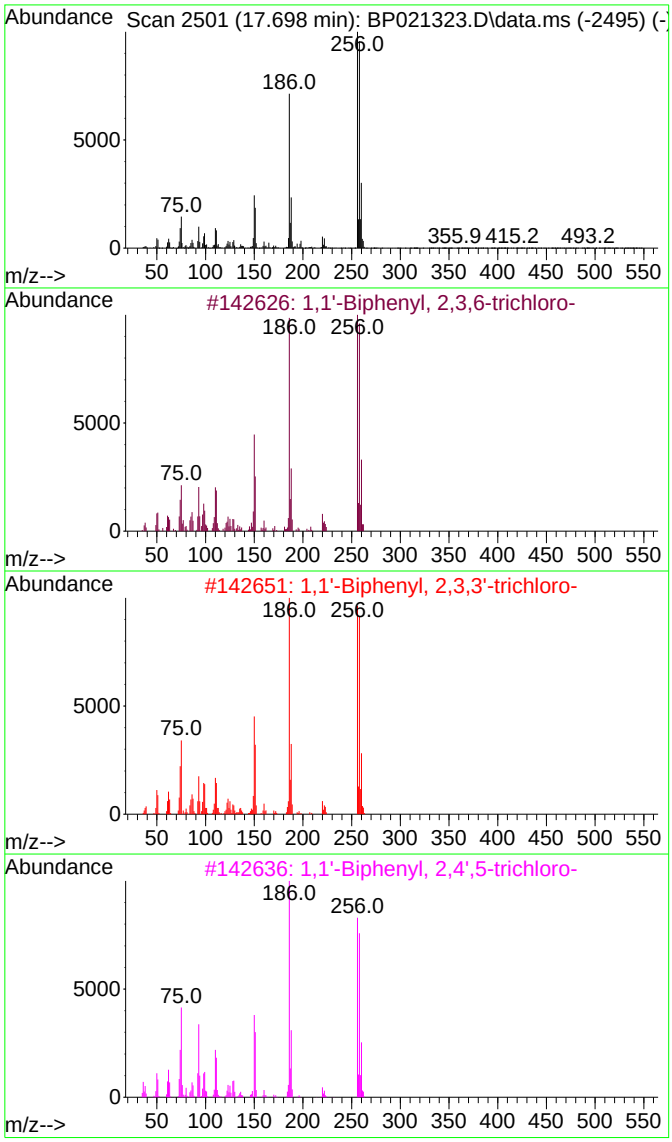
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

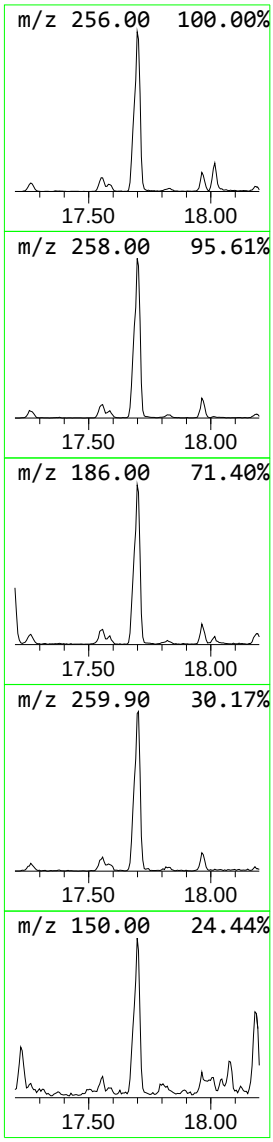
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1,1'-Biphenyl, 2,3,6-trichl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.698	7.24 ng/ul	795309	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	99
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
4	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
5	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

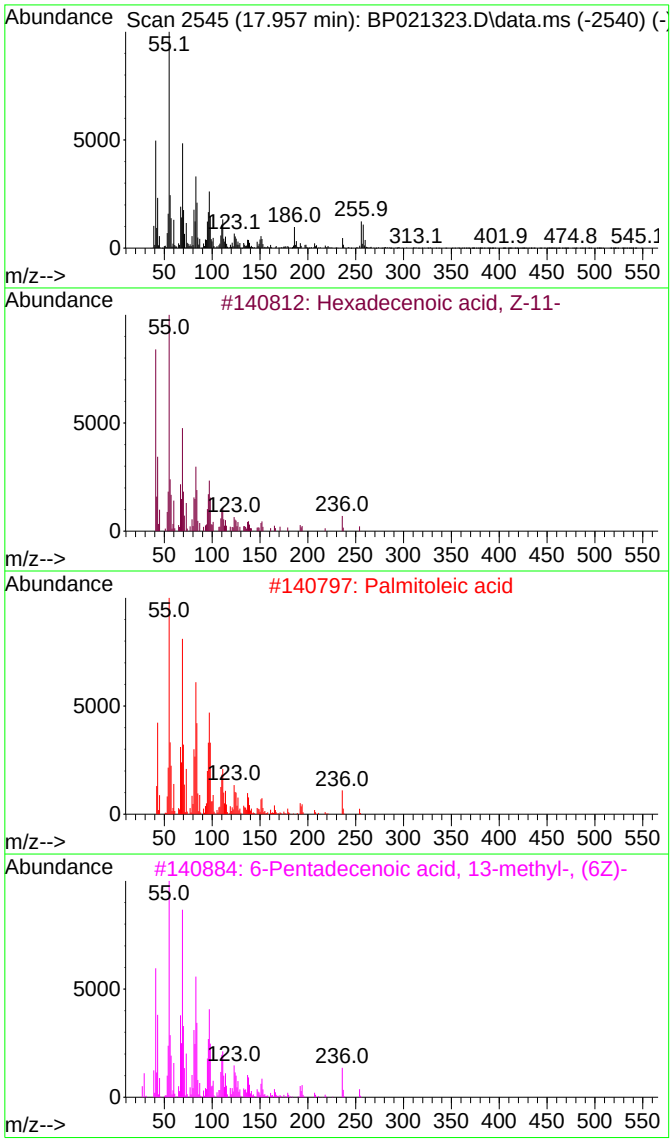
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

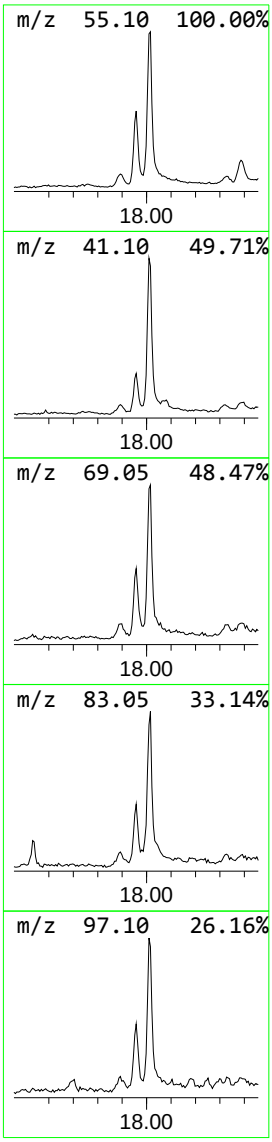
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexadecenoic acid, Z-11- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.957	4.13 ng/ul	453191	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	98
2	Palmitoleic acid	254	C16H30O2	000373-49-9	96
3	6-Pentadecenoic acid, 13-methyl-...	254	C16H30O2	682751-34-4	89
4	cis-Vaccenic acid	282	C18H34O2	000506-17-2	60
5	cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	53





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Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

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Quant Title : SVOA CALIBRATION

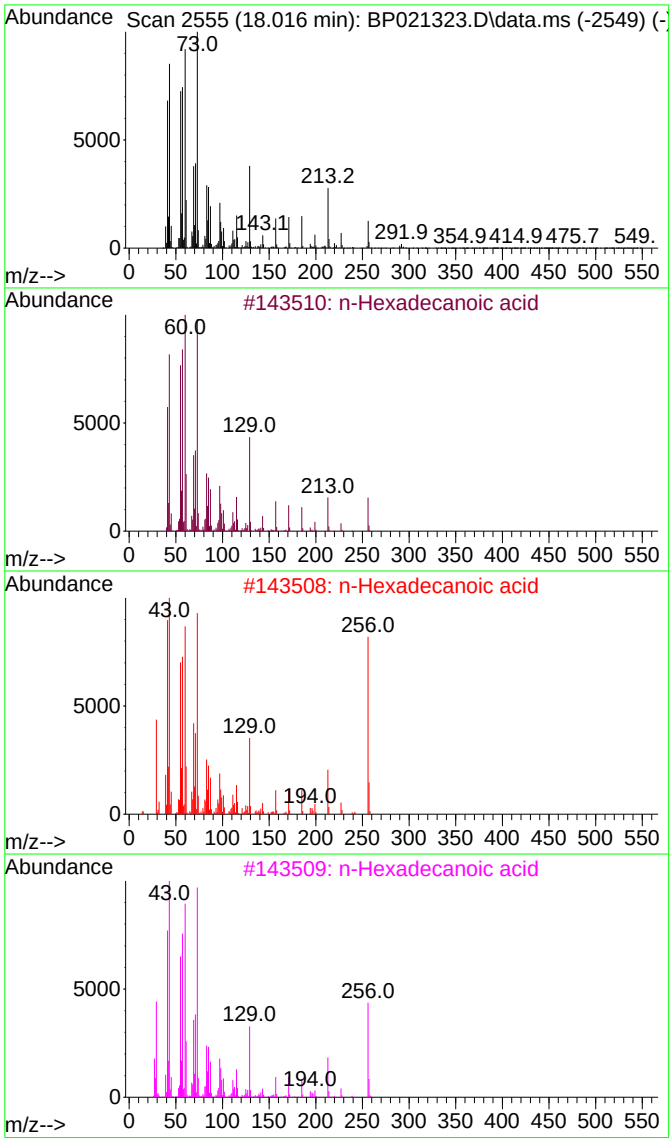
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

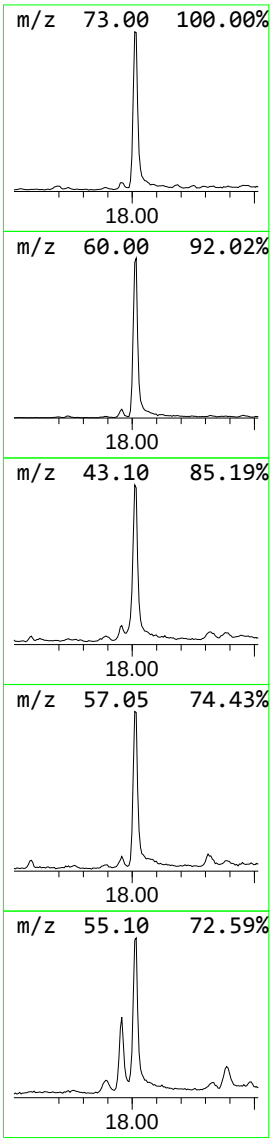
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.016	15.64 ng/ul	1717300	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	Octadecanoic acid	284	C18H36O2	000057-11-4	91





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Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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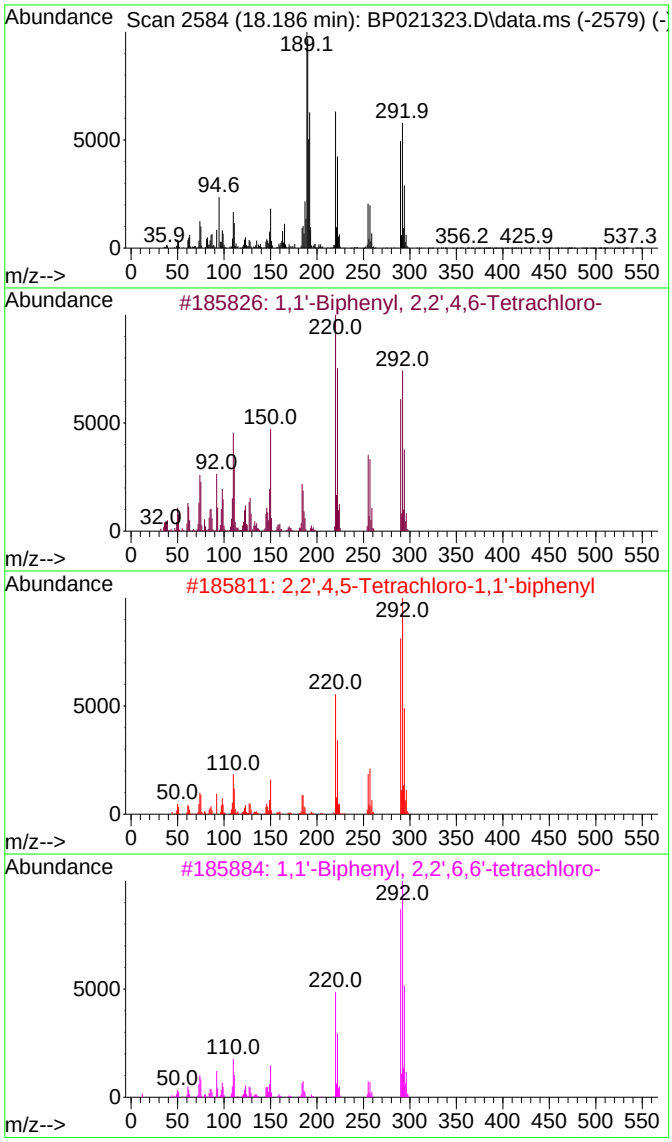
Quant Title : SVOA CALIBRATION

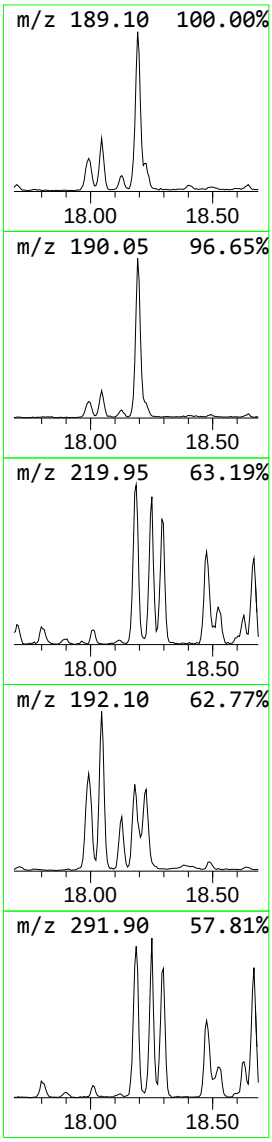
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1,1'-Biphenyl, 2,2',4,6-Tet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.186	7.57 ng/ul	831069	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	96
2	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	96
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	96
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	96
5	1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	96





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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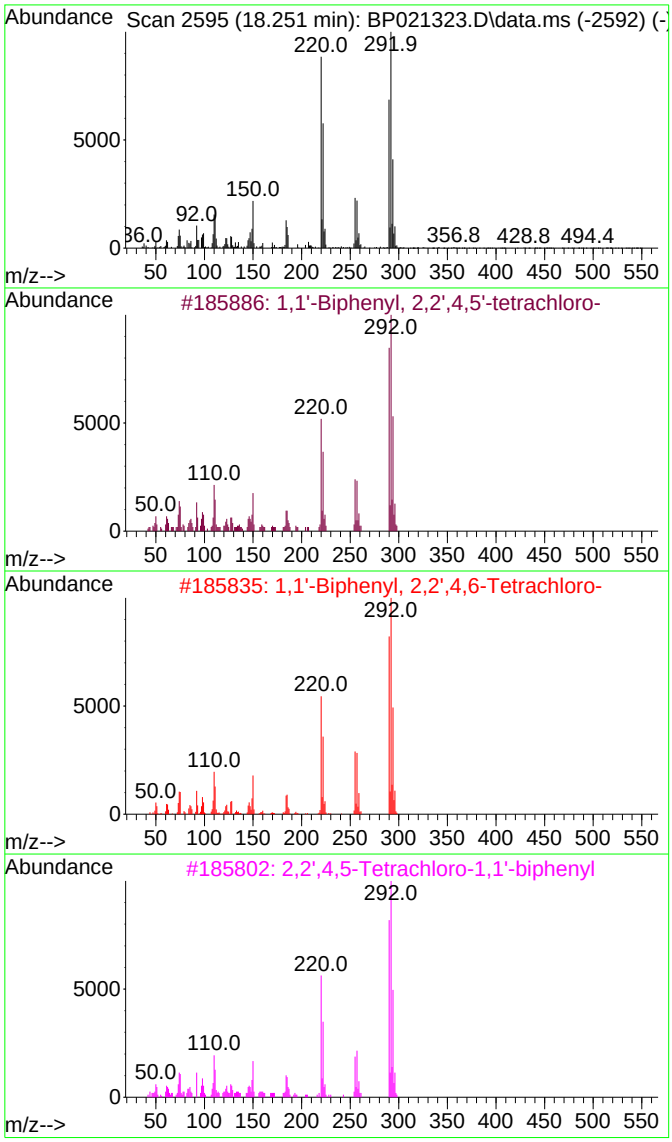
Quant Title : SVOA CALIBRATION

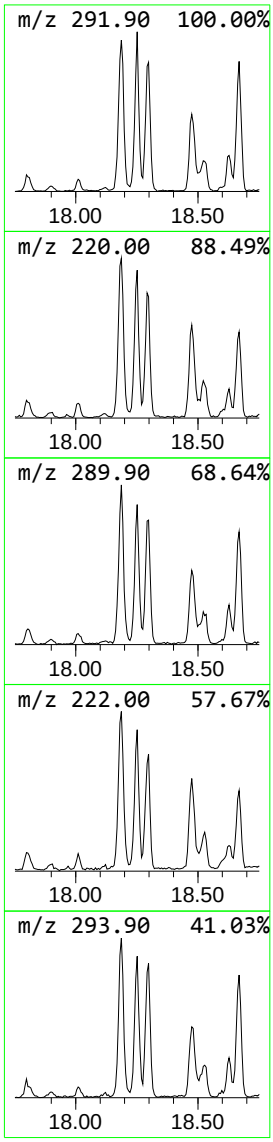
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,1'-Biphenyl, 2,2',4,5'-te... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.251	2.49 ng/ul	273900	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
2	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99
3	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
4	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	99
5	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99





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Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

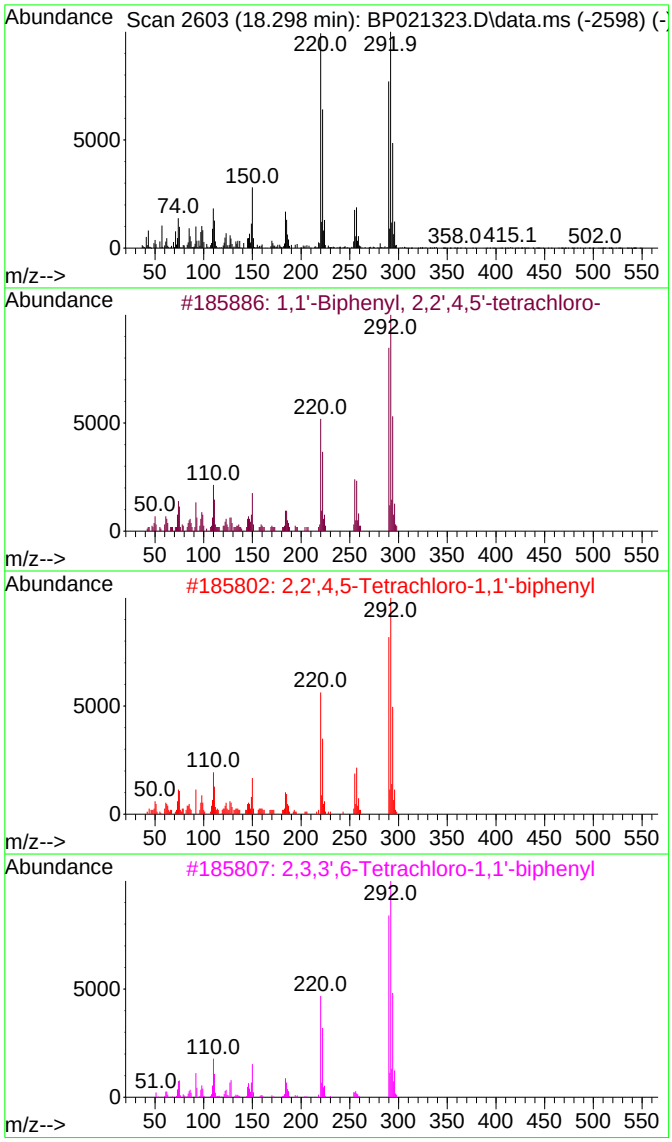
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

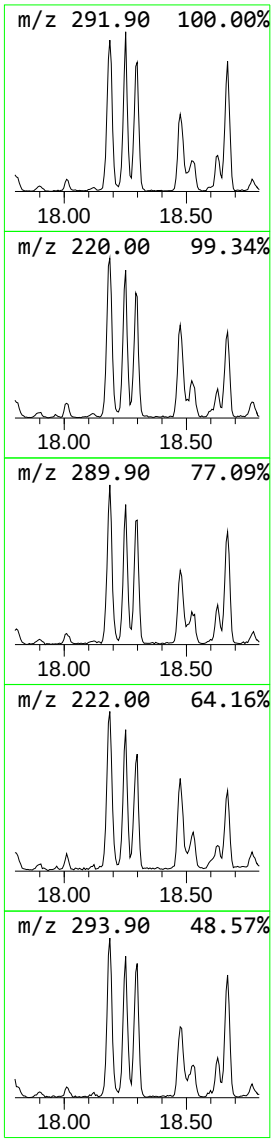
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.298	2.58 ng/ul	283338	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
2	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
3	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99
4	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
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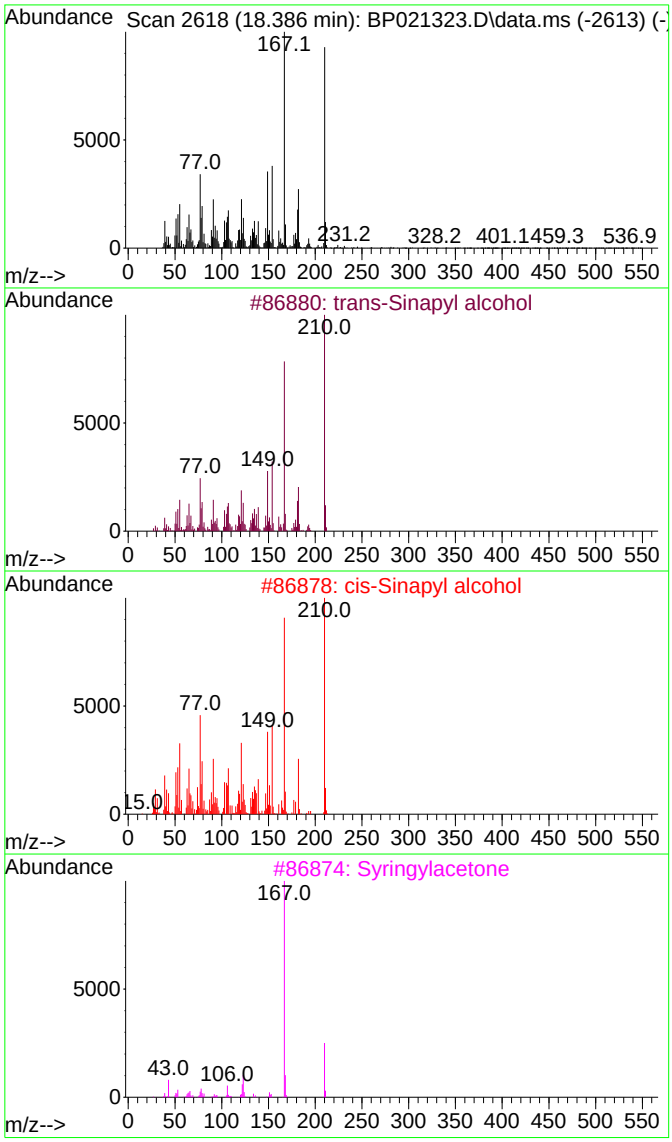
Quant Title : SVOA CALIBRATION

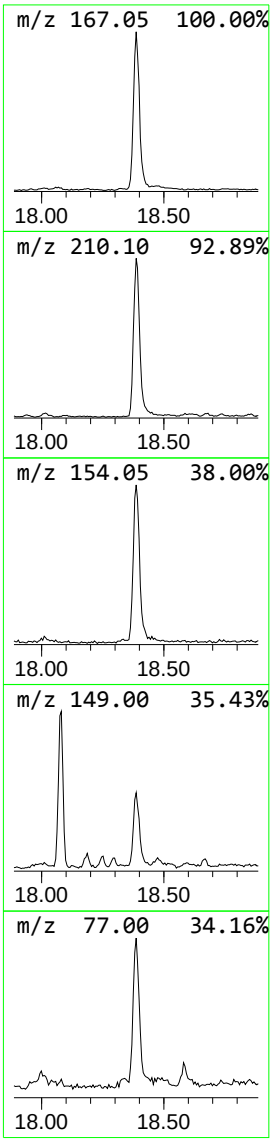
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 trans-Sinapyl alcohol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.386	9.87 ng/ul	1084240	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-Sinapyl alcohol	210	C11H14O4	020675-96-1	99
2	cis-Sinapyl alcohol	210	C11H14O4	104330-63-4	94
3	Syringylacetone	210	C11H14O4	019037-58-2	64
4	4-Ethoxy-2,5-dimethoxybenzaldehyde	210	C11H14O4	1000121-30-2	58
5	2-Pentanone, 1-(2,4,6-trihydroxy...	210	C11H14O4	1000116-22-3	52





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Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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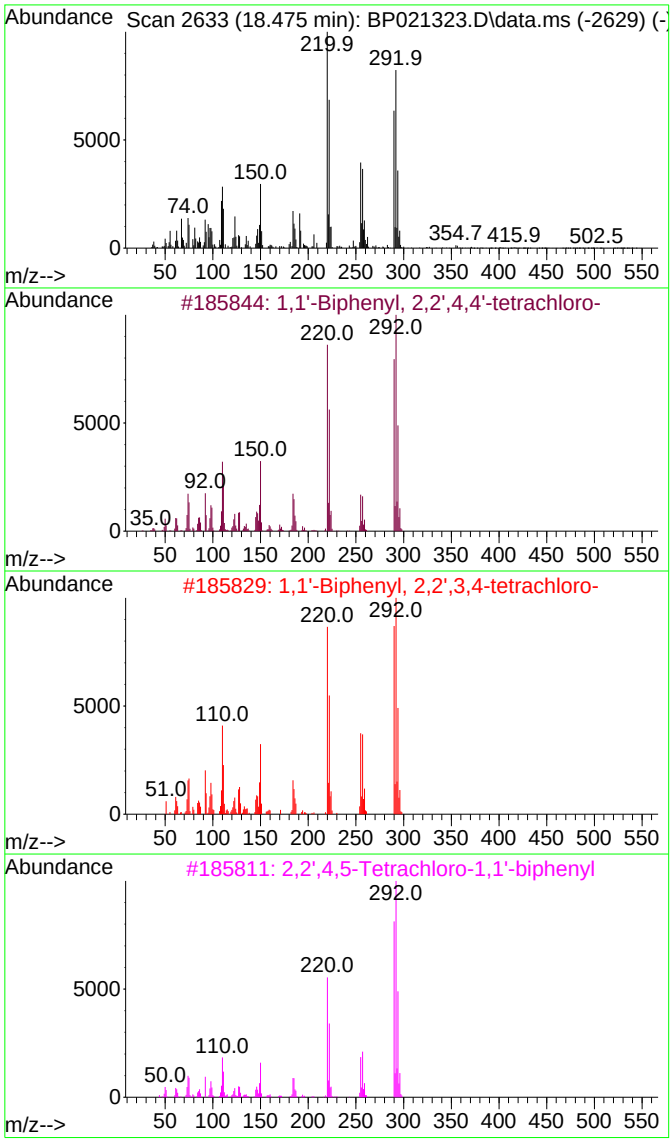
Quant Title : SVOA CALIBRATION

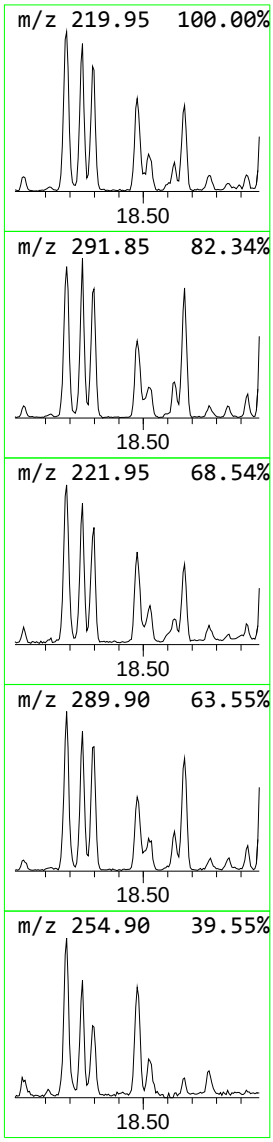
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.475	2.22 ng/ul	243418	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
3	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99





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Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

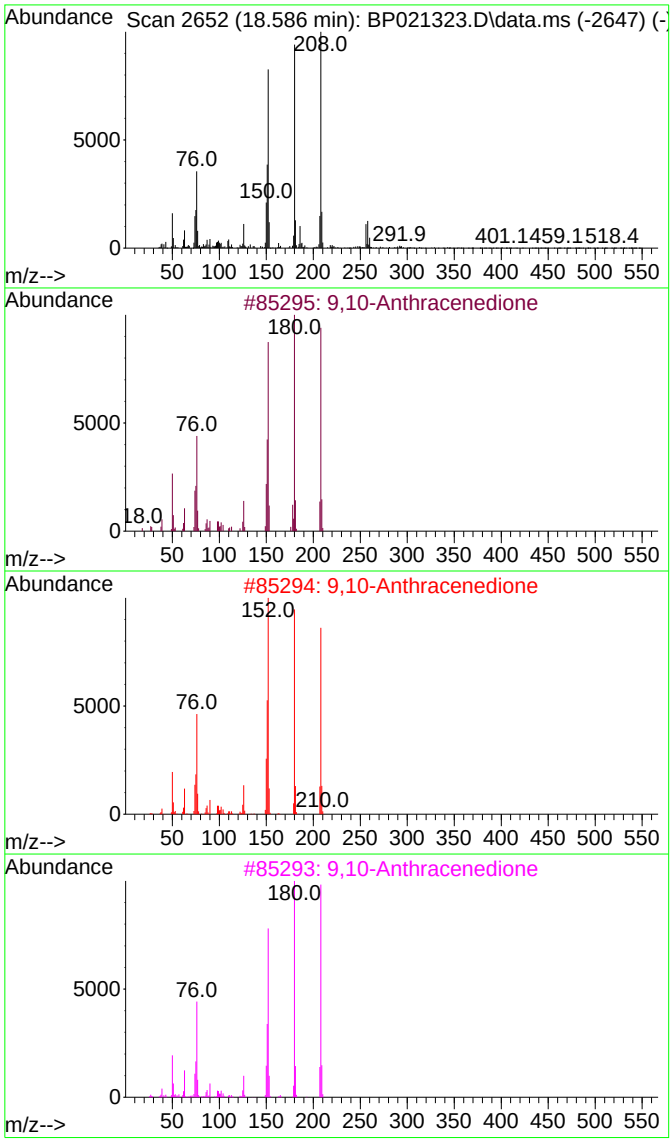
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

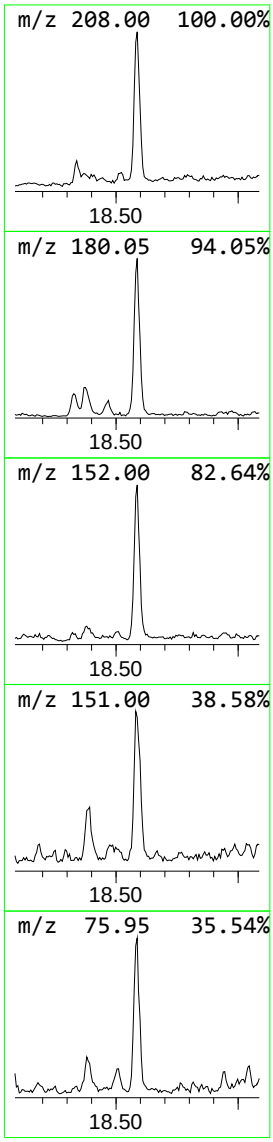
****

Peak Number 12 9,10-Anthracenedione Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.586	3.61 ng/ul	396445	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	93
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	86
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	83





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Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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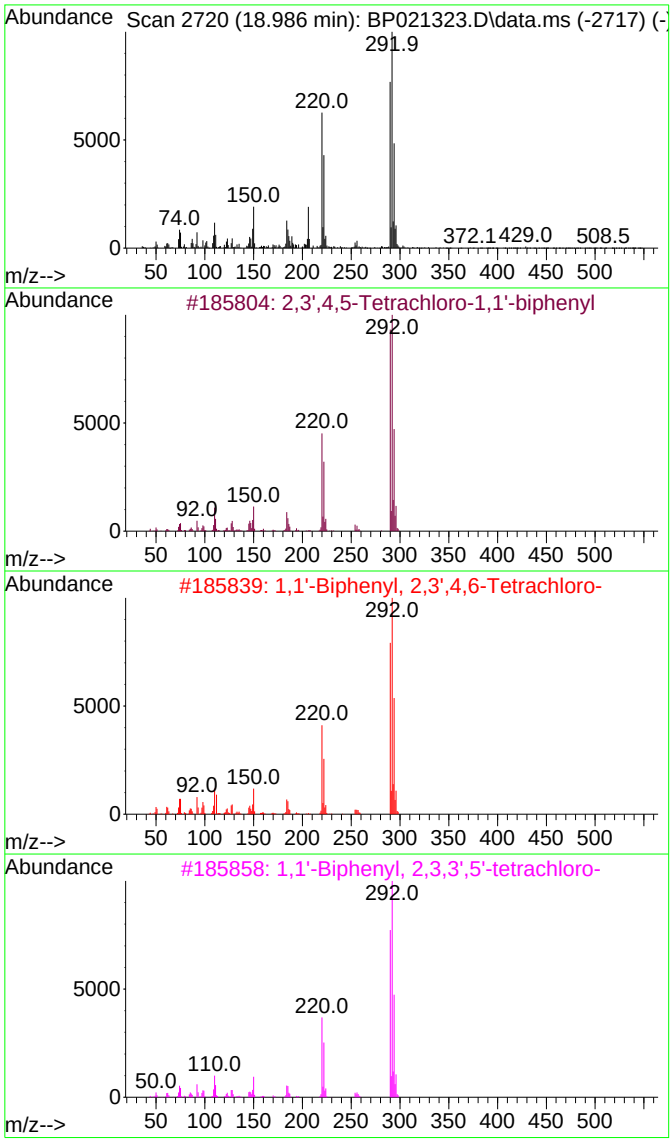
Quant Title : SVOA CALIBRATION

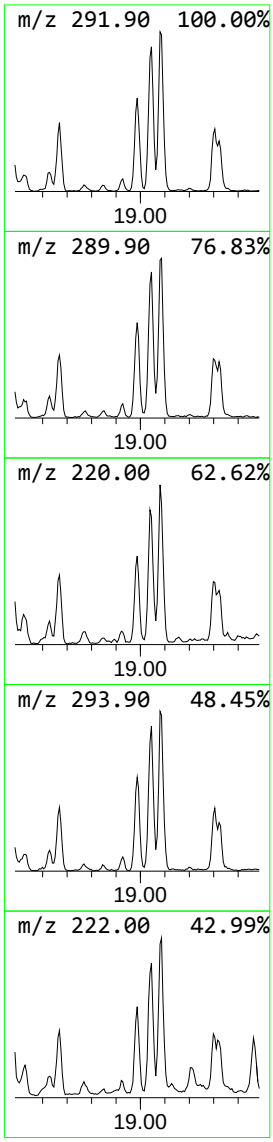
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 2,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.986	2.31 ng/ul	254011	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
2	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99
3	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
4	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
5	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	99





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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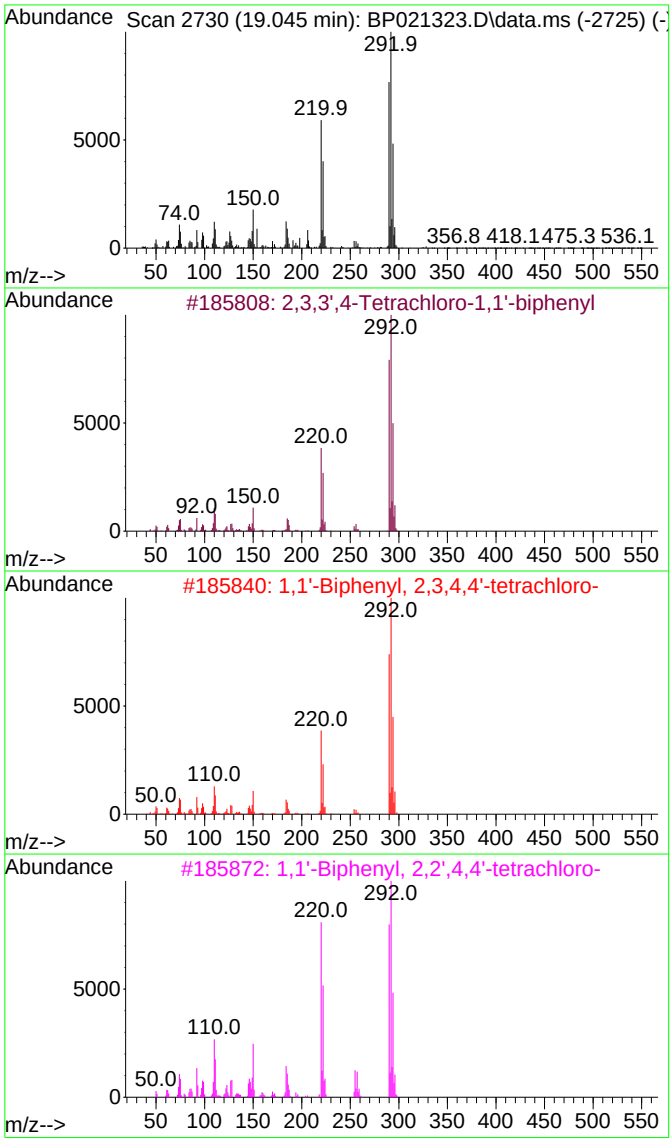
Quant Title : SVOA CALIBRATION

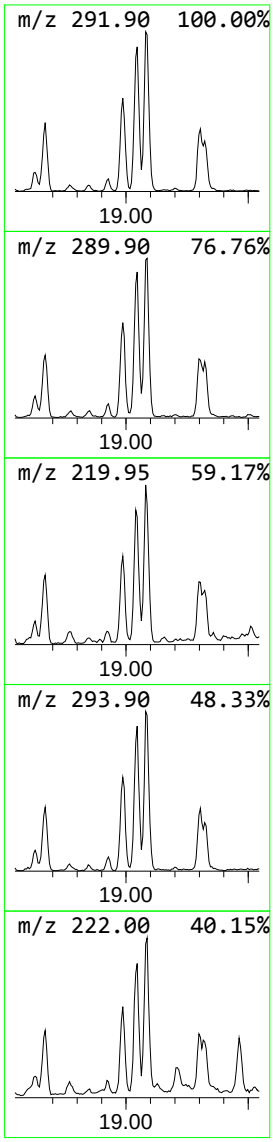
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 2,3,3',4-Tetrachloro-1,1'-b... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.045	4.39 ng/ul	482693	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',4-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-24-2	99
2	1,1'-Biphenyl, 2,3,4,4'-tetrachl...	290	C12H6Cl4	033025-41-1	99
3	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
4	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99
5	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	99





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ALS Vial : 5 Sample Multiplier: 1

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ClientSampleId :
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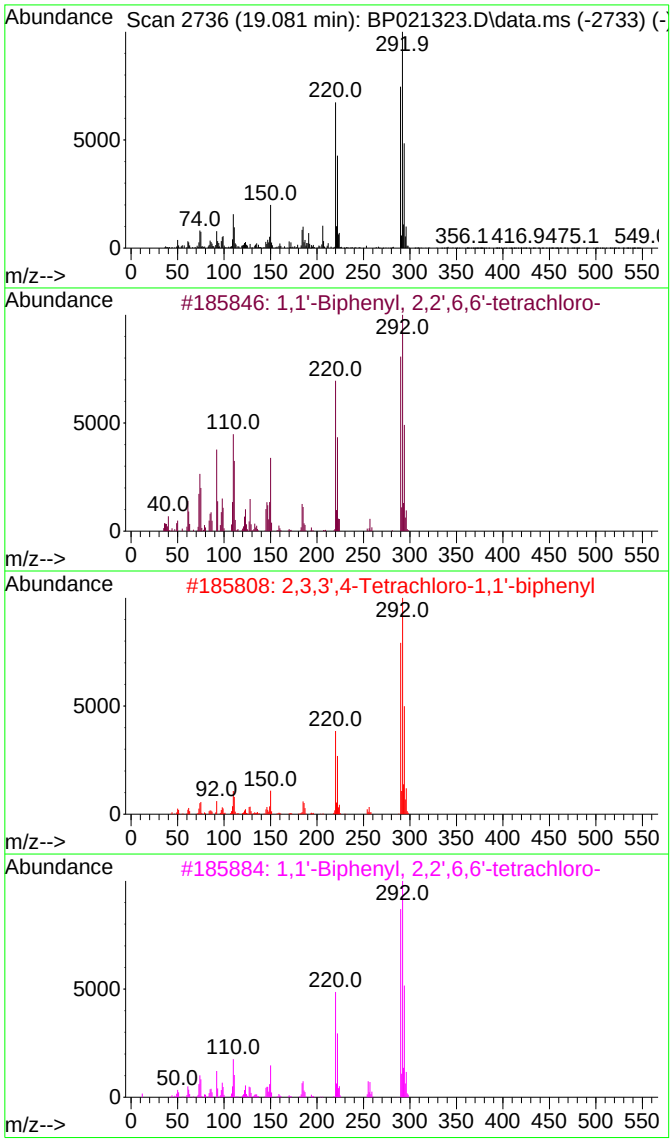
Quant Title : SVOA CALIBRATION

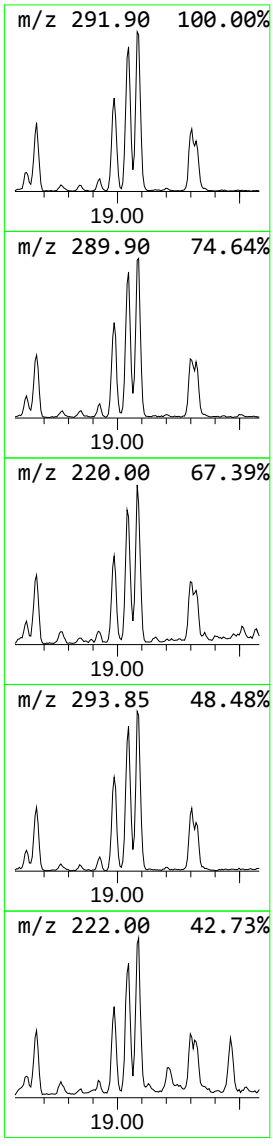
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 1,1'-Biphenyl, 2,2',6,6'-te... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.081	3.52 ng/ul	386171	Phenanthrene-d10	17.087

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	95
2	2,3,3',4-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-24-2	95
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	95
4	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	95
5	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	95





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Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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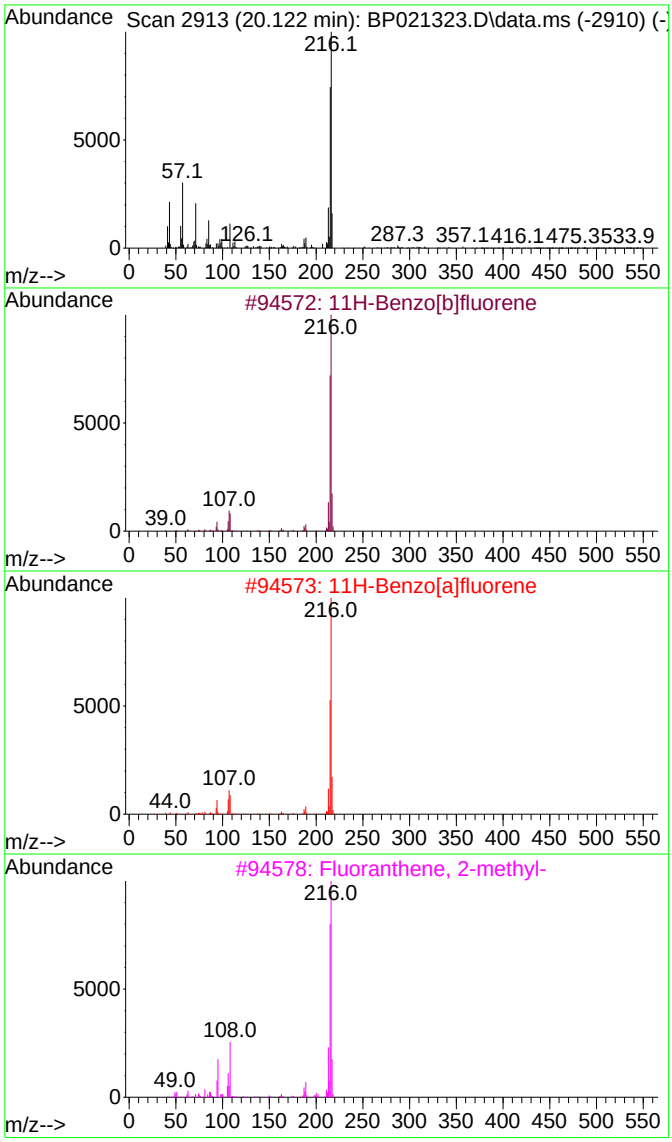
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

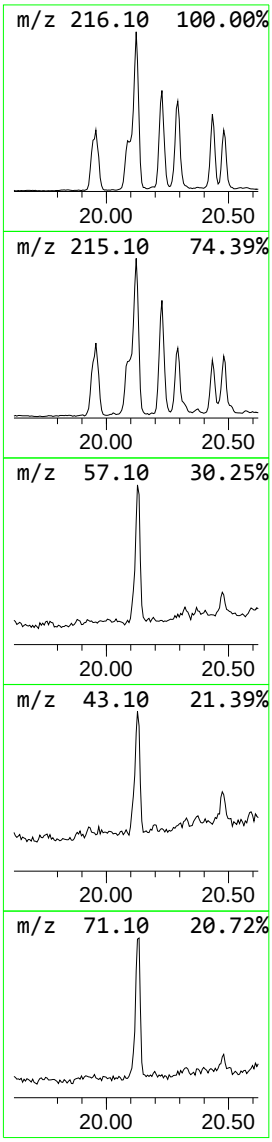
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 11H-Benzo[b]fluorene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.122	2.16 ng/ul	660049	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76





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Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

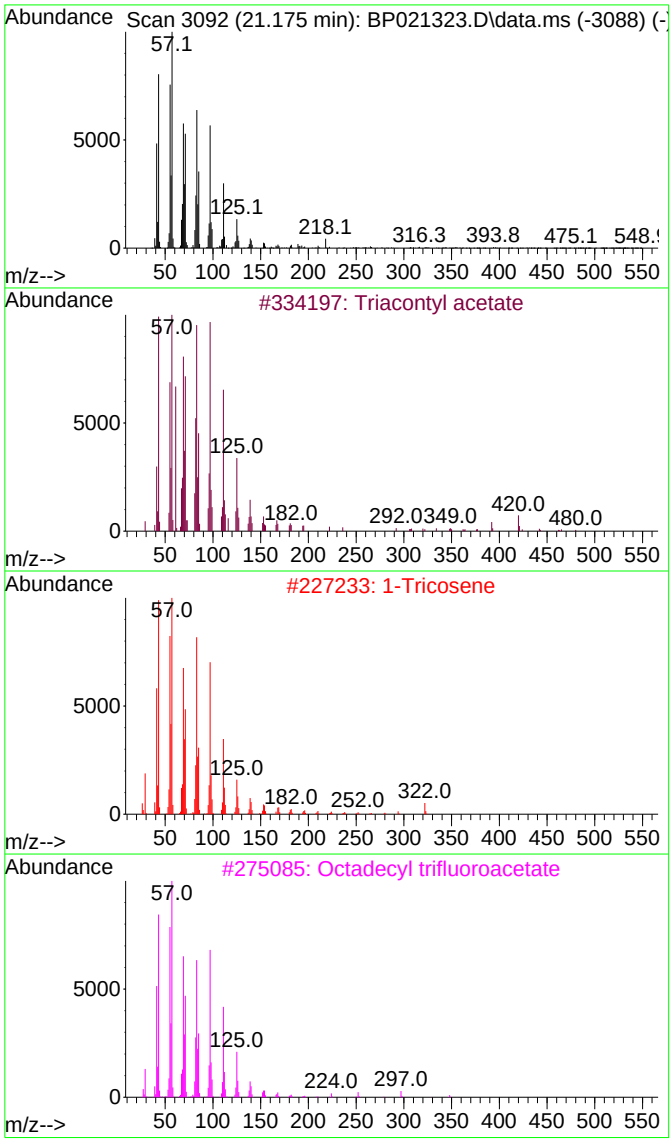
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

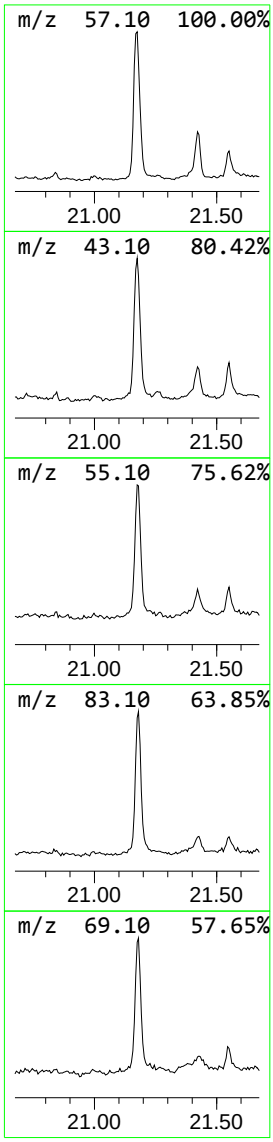
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Triacontyl acetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.174	7.65 ng/ul	2336150	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontyl acetate	480	C32H64O2	041755-58-2	96
2	1-Tricosene	322	C23H46	018835-32-0	95
3	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91





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Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

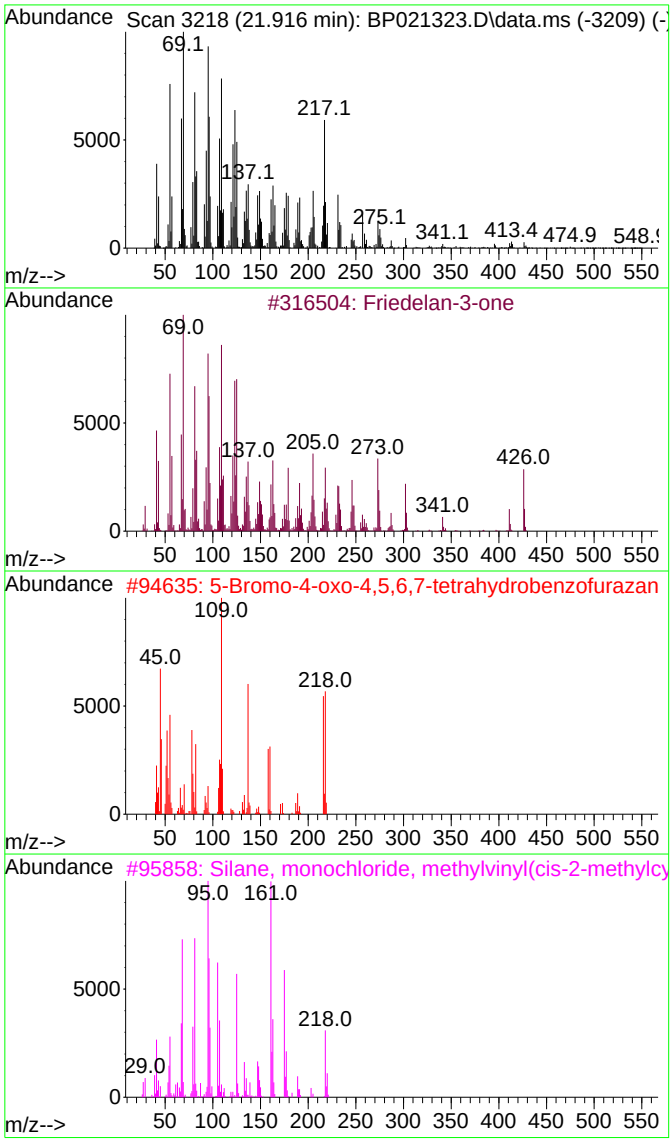
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

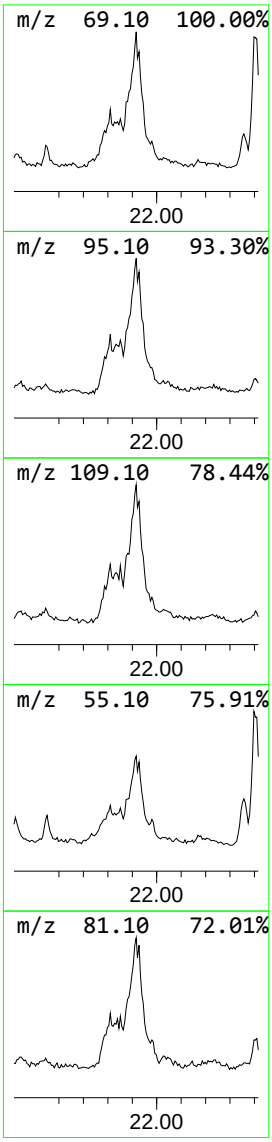
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.916	12.46 ng/ul	3804180	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Friedelan-3-one	426	C30H50O	000559-74-0	89
2	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	70
3	Silane, monochloride, methylviny...	218	C10H19ClOSi	1000421-06-3	70
4	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	70
5	2,4a,8,8-Tetramethyldecahydrocyc...	206	C15H26	074022-04-1	59





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Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

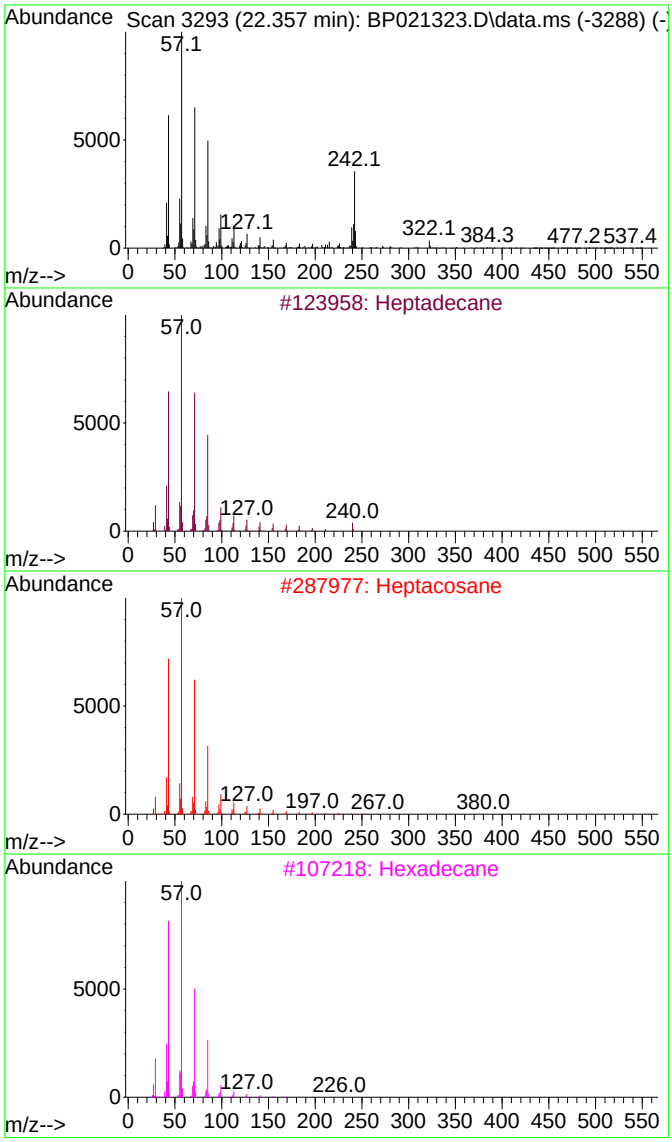
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

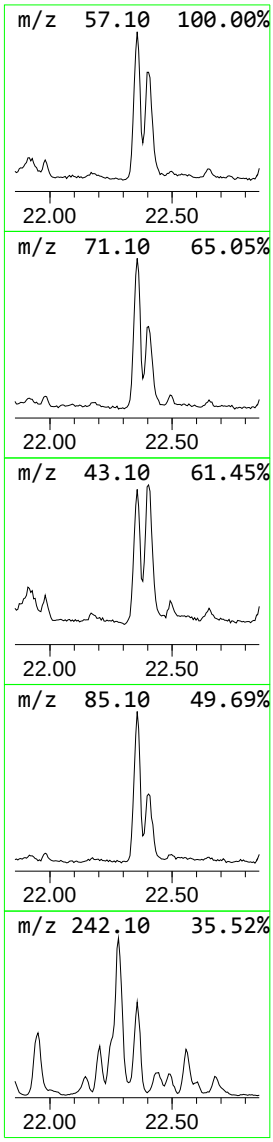
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.357	4.15 ng/ul	1267360	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	98
2	Heptacosane	380	C27H56	000593-49-7	96
3	Hexadecane	226	C16H34	000544-76-3	92
4	Heptadecane	240	C17H36	000629-78-7	91
5	Hexacosane	366	C26H54	000630-01-3	91





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Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

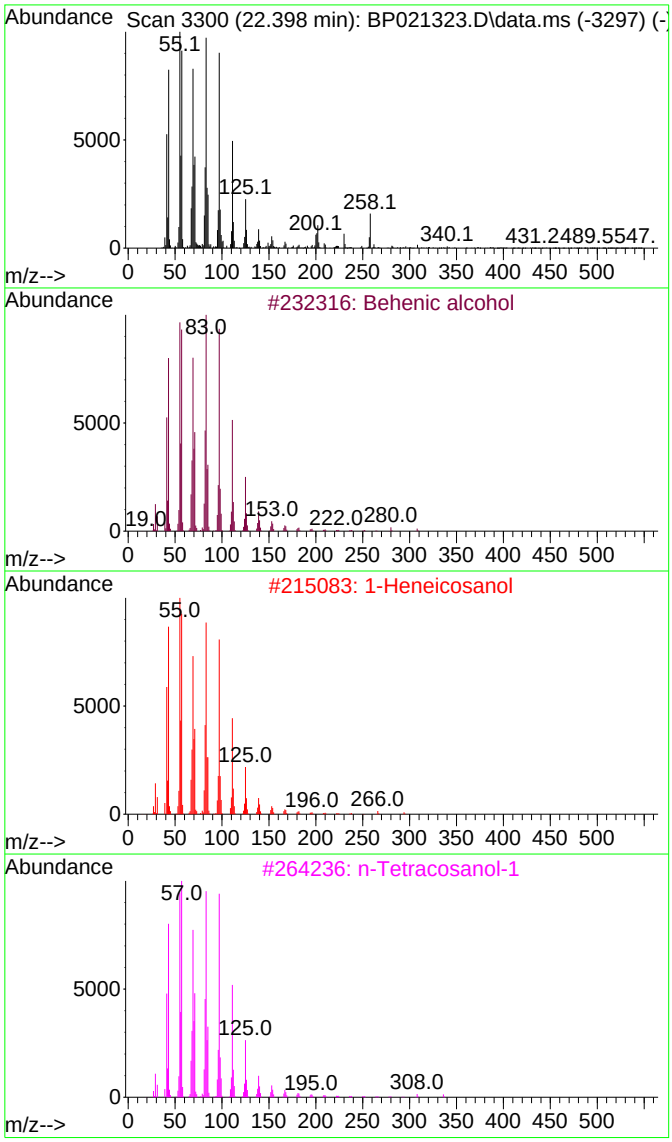
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

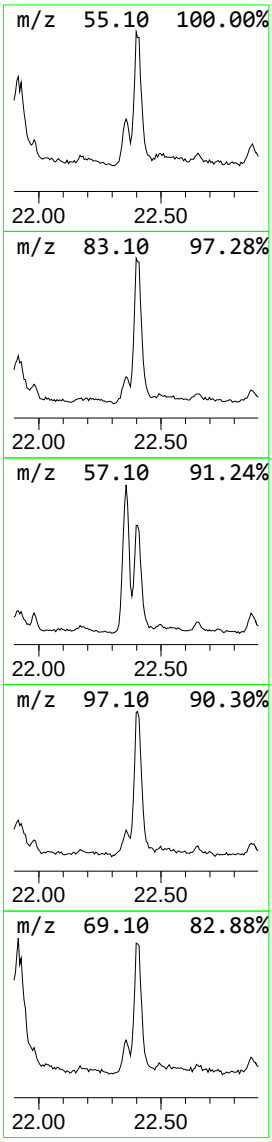
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Behenic alcohol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.398	7.21 ng/ul	2202620	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	1-Hexacosene	364	C26H52	018835-33-1	93
5	n-Heptadecanol-1	256	C17H36O	001454-85-9	93





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Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

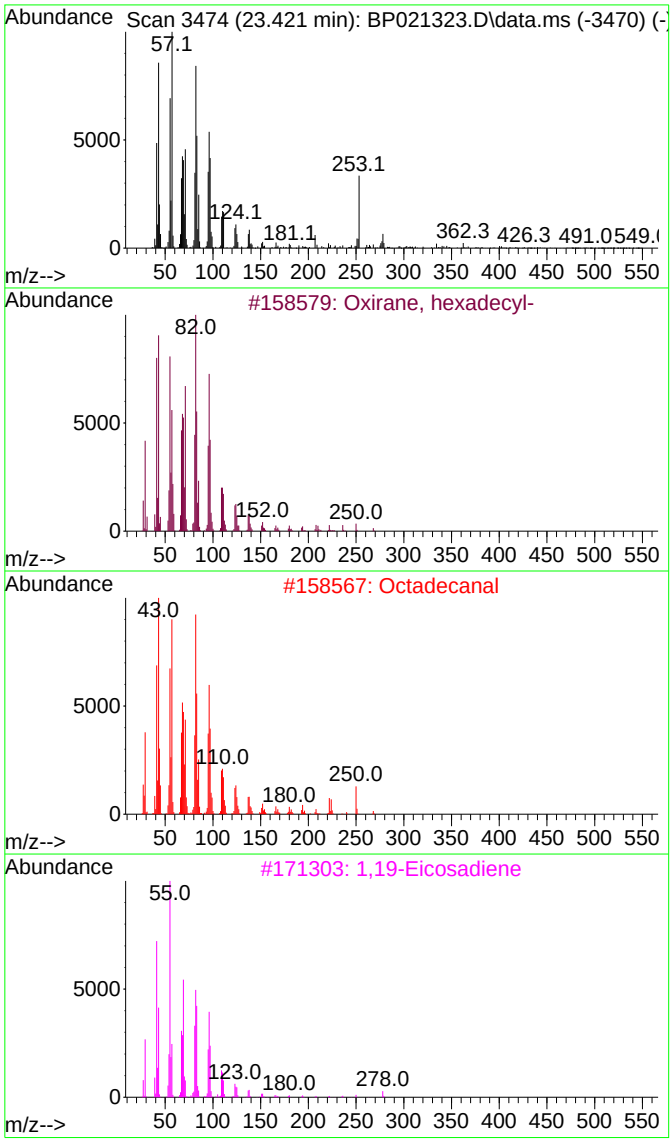
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

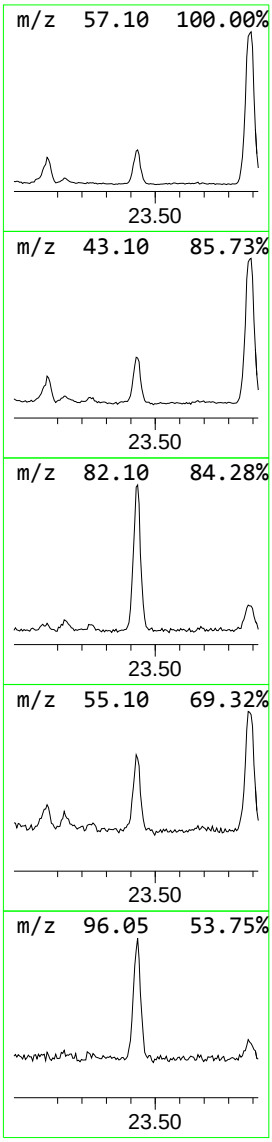
****

Peak Number 21 Oxirane, hexadecyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.422	8.04 ng/ul	1230860	Perylene-d12	24.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	95
2	Octadecanal	268	C18H36O	000638-66-4	95
3	1,19-Eicosadiene	278	C20H38	014811-95-1	94
4	Octacosanal	408	C28H56O	022725-64-0	94
5	Hexacosanal	380	C26H52O	026627-85-0	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

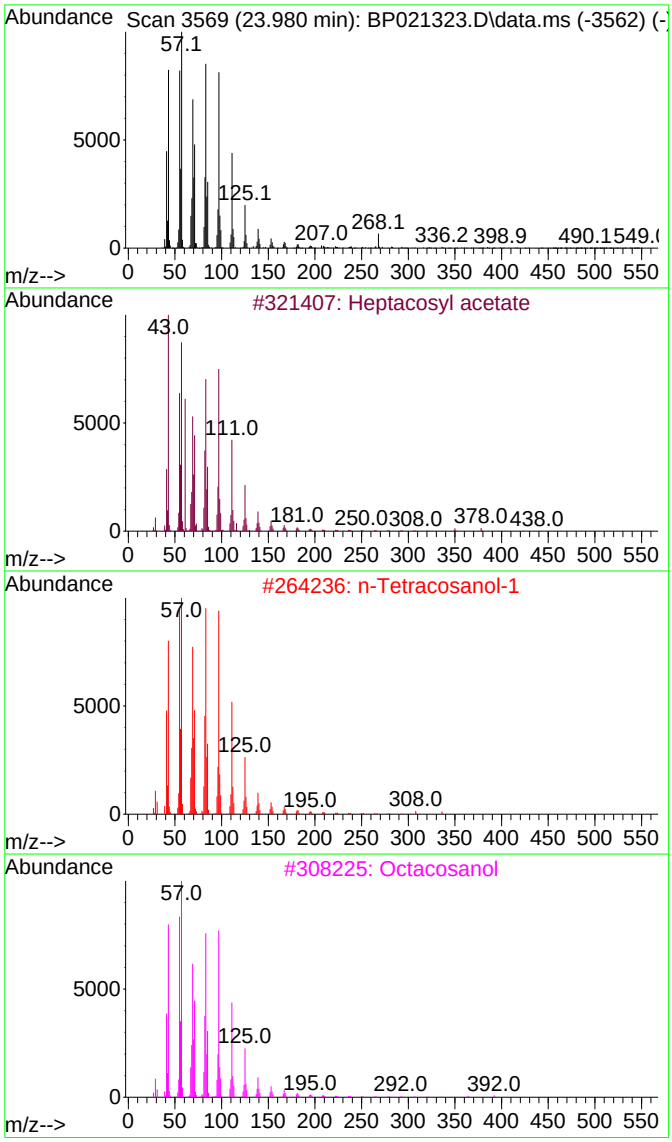
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

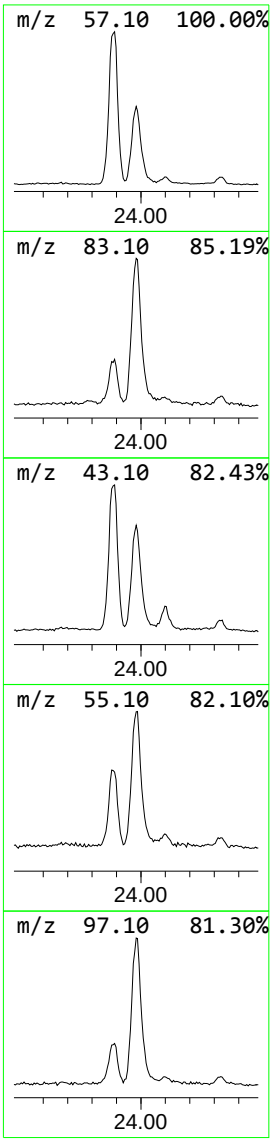
****

Peak Number 22 Heptacosyl acetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.980	20.12 ng/ul	3080780	Perylene-d12	24.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosyl acetate	438	C29H58O2	1000351-78-2	97
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3	Octacosanol	410	C28H58O	000557-61-9	94
4	1-Heptacosanol	396	C27H56O	002004-39-9	94
5	1-Hexacosanol	382	C26H54O	000506-52-5	94





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Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

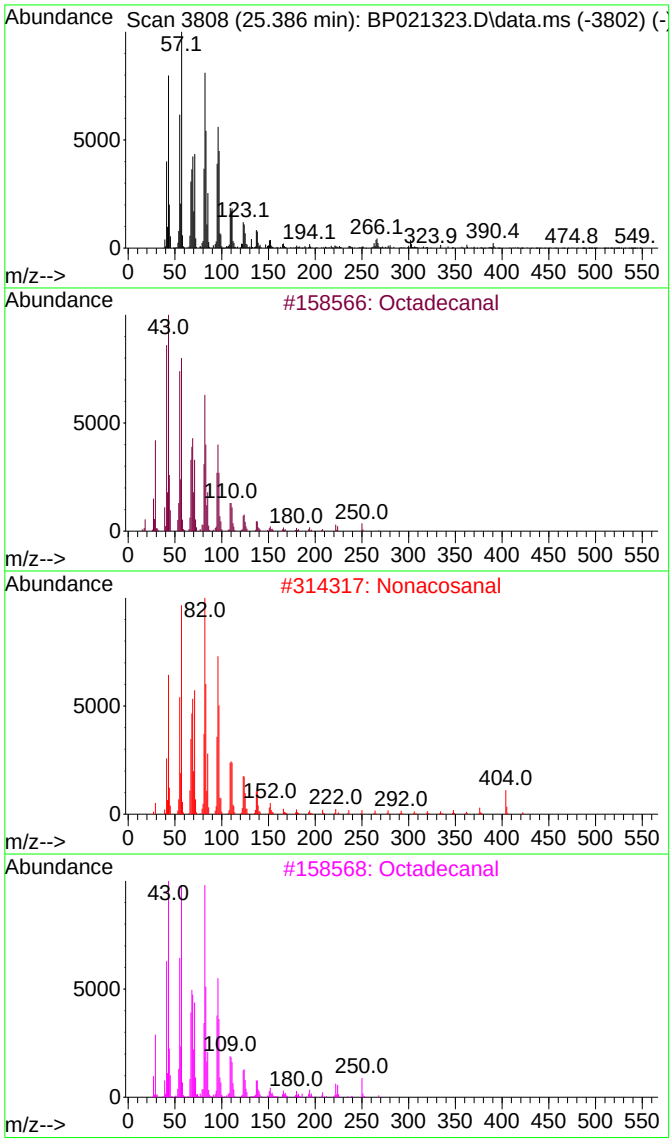
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

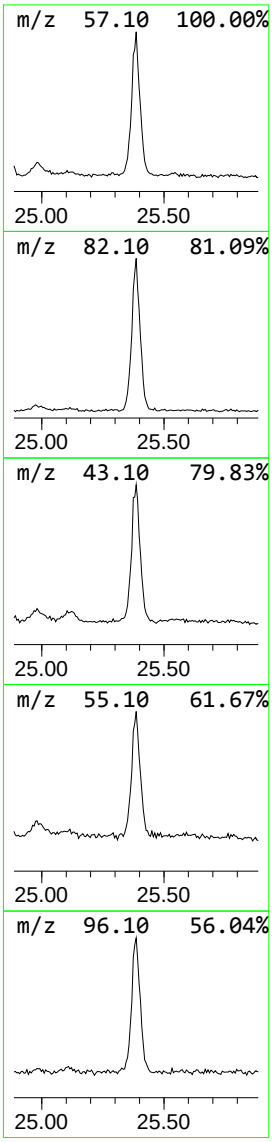
****

Peak Number 23 Octadecanal Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.386	14.77 ng/ul	2262070	Perylene-d12	24.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanal	268	C18H36O	000638-66-4	94
2	Nonacosanal	422	C29H58O	072934-04-4	93
3	Octadecanal	268	C18H36O	000638-66-4	91
4	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
5	Octadecanal	268	C18H36O	000638-66-4	91





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Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

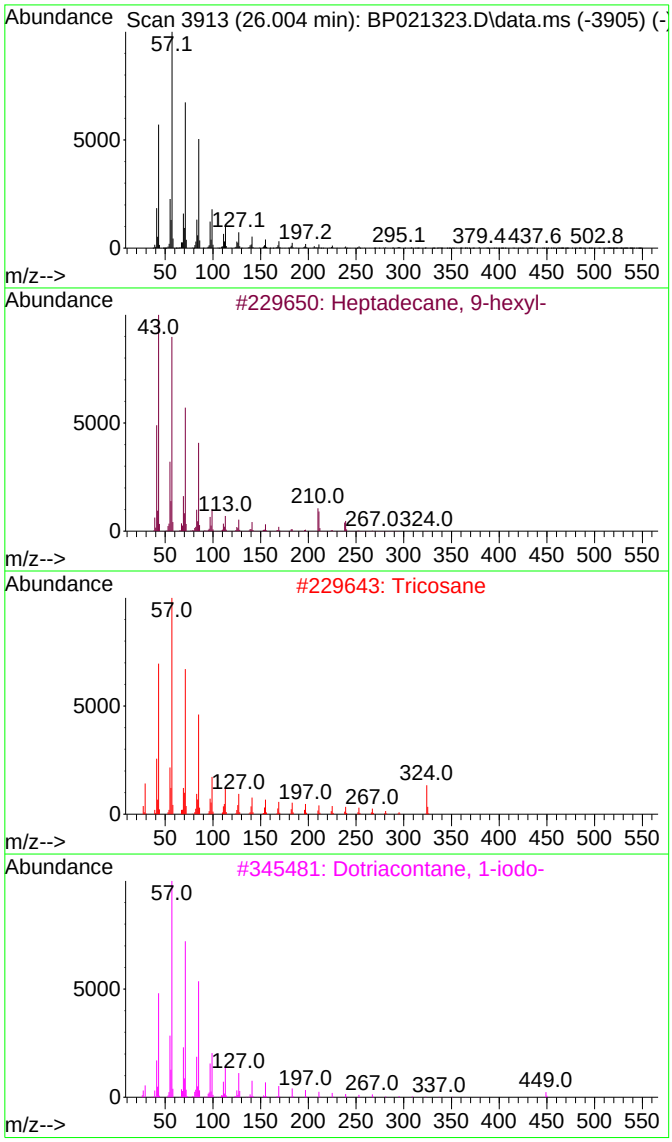
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

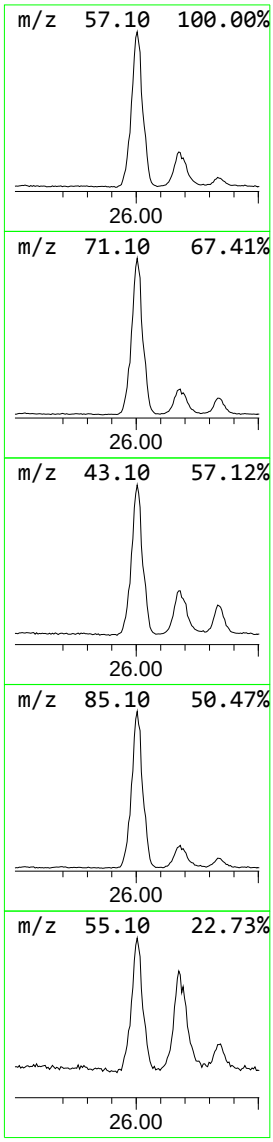
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.004	28.77 ng/ul	4404690	Perylene-d12	24.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2	Tricosane	324	C23H48	000638-67-5	93
3	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
4	Tetratetracontane	619	C44H90	007098-22-8	91
5	Triacontane	422	C30H62	000638-68-6	91





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Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

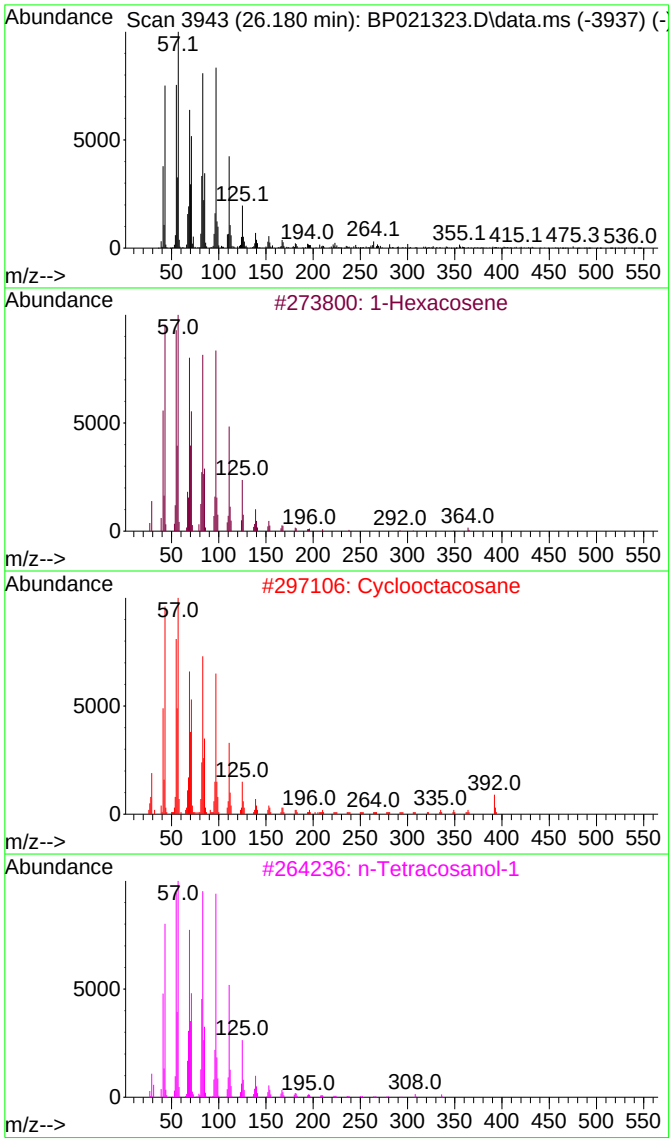
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

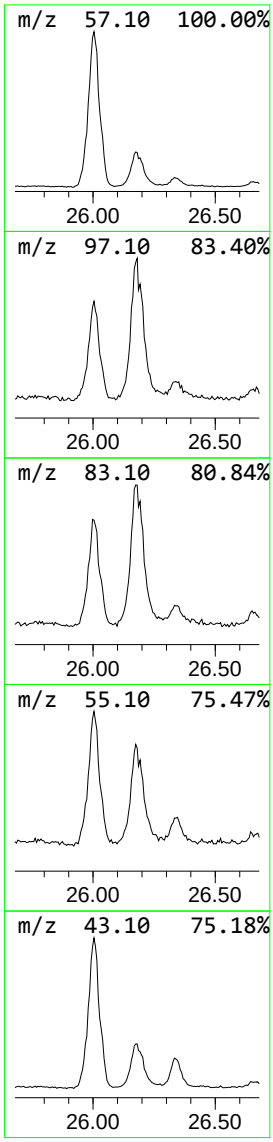
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.180	13.36 ng/ul	2045730	Perylene-d12	24.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	94
2	Cyclooctacosane	392	C28H56	000297-24-5	93
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
5	1-Heptacosanol	396	C27H56O	002004-39-9	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080124\
Data File : BP021323.D
Acq On : 02 Aug 2024 16:18
Operator : CG/JU
Sample : P3265-06
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4D39

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
Benzeneethanol,...	13.499	2.2	ng/ul	179823	3	14.263	1630210	20.0
unknown-01	14.487	4.4	ng/ul	357824	3	14.263	1630210	20.0
(E)-4-(3-Hydrox...	16.492	15.6	ng/ul	1709920	4	17.087	2196680	20.0
1,1'-Biphenyl, ...	17.698	7.2	ng/ul	795309	4	17.087	2196680	20.0
Hexadecenoic ac...	17.957	4.1	ng/ul	453191	4	17.087	2196680	20.0
n-Hexadecanoic ...	18.016	15.6	ng/ul	1717300	4	17.087	2196680	20.0
1,1'-Biphenyl, ...	18.186	7.6	ng/ul	831069	4	17.087	2196680	20.0
1,1'-Biphenyl, ...	18.251	2.5	ng/ul	273900	4	17.087	2196680	20.0
2,2',4,5-Tetrac...	18.298	2.6	ng/ul	283338	4	17.087	2196680	20.0
trans-Sinapyl a...	18.386	9.9	ng/ul	1084240	4	17.087	2196680	20.0
1,1'-Biphenyl, ...	18.475	2.2	ng/ul	243418	4	17.087	2196680	20.0
9,10-Anthracene...	18.586	3.6	ng/ul	396445	4	17.087	2196680	20.0
2,3',4,5-Tetrac...	18.986	2.3	ng/ul	254011	4	17.087	2196680	20.0
2,3,3',4-Tetrac...	19.045	4.4	ng/ul	482693	4	17.087	2196680	20.0
1,1'-Biphenyl, ...	19.081	3.5	ng/ul	386171	4	17.087	2196680	20.0
11H-Benzo[b]flu...	20.122	2.2	ng/ul	660049	5	21.533	6108110	20.0
Triacontyl acetate	21.174	7.7	ng/ul	2336150	5	21.533	6108110	20.0
5-Bromo-4-oxo-4...	21.916	12.5	ng/ul	3804180	5	21.533	6108110	20.0
(DEL) Alkane: S...	22.357	4.2	ng/ul	1267360	5	21.533	6108110	20.0
Behenic alcohol	22.398	7.2	ng/ul	2202620	5	21.533	6108110	20.0
Oxirane, hexade...	23.422	8.0	ng/ul	1230860	6	24.827	3062400	20.0
Heptacosyl acetate	23.980	20.1	ng/ul	3080780	6	24.827	3062400	20.0
Octadecanal	25.386	14.8	ng/ul	2262070	6	24.827	3062400	20.0
(DEL) Alkane: S...	26.004	28.8	ng/ul	4404690	6	24.827	3062400	20.0
(DEL) Alkane: S...	26.180	13.4	ng/ul	2045730	6	24.827	3062400	20.0