LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A4CG7

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: BP021158.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	30	33	39	rVB	32611	45667	0.48%	0.056%
2	3.469	79	82	96	rVB	114819	176947	1.87%	0.216%
3	3.622	105	108	118	rBV	24945	44420	0.47%	0.054%
4	3.899	149	155	161	rVB2	15300	29039	0.31%	0.035%
5	4.269	214	218	223	rBV7	7297	10616	0.11%	0.013%
6	4.804	305	309	316	rBV	12622	21544	0.23%	0.026%
7	6.551	599	606	611	rBV3	7459	13479	0.14%	0.016%
8	6.757	634	641	649	rBV7	11988	35534	0.38%	0.043%
9	6.875	655	661	677	rBV	467992	838041	8.86%	1.023%
10	6.998	677	682	695	rVB	402553	631020	6.67%	0.771%
11	7.204	711	717	724	rBV	568335	886593	9.38%	1.083%
12	7.275	726	729	738	rVB	37224	77722	0.82%	0.095%
13	7.651	786	793	803	rBV	779338	1209456	12.79%	1.477%
14	7.863	821	829	836	rBV6	10985	28297	0.30%	0.035%
15	8.387	911	918	934	rBV	558739	1053030	11.14%	1.286%
16	8.810	983	990	1004	rBV	436694	816837	8.64%	0.997%
17	8.963	1010	1016	1026	rVB5	20157	41502	0.44%	0.051%
18	9.369	1080	1085	1091	rBV2	15668	27148	0.29%	0.033%
19	9.522	1101	1111	1127	rBV	372413	731088	7.73%	0.893%
20	9.810	1153	1160	1170	rBV7	21350	63465	0.67%	0.078%
21	10.087	1199	1207	1224	rBV	501683	1085879	11.48%	1.326%
22	10.422	1249	1264	1277	rBV	944107	1693449	17.91%	2.068%
23	10.587	1286	1292	1309	rVB	63604	156524	1.66%	0.191%
24	10.828	1326	1333	1338	rBV	5422	15697	0.17%	0.019%
25	10.963	1351	1356	1363	rBV4	19222	32991	0.35%	0.040%
26	11.057	1363	1372	1381	rBV	276493	834762	8.83%	1.019%
27	11.181	1387	1393	1402	rVB2	146498	323050	3.42%	0.394%
28	11.251	1402	1405	1412	rVB5	17562	33360	0.35%	0.041%
29	11.822	1495	1502	1509	rBV7	15895	36936	0.39%	0.045%
30	12.016	1532	1535	1543	rVB3	11810	22393	0.24%	0.027%
31	12.128	1548	1554	1559	rBV	33447	63204	0.67%	0.077%
32	12.745	1652	1659	1662	rBV9	8426	20787	0.22%	0.025%
33	12.792	1665	1667	1675	rVV9	9296	21552	0.23%	0.026%
34	12.875	1677	1681	1688	rVB10	10624	23764	0.25%	0.029%
35	13.069	1706	1714	1721	rBV2	97221	199155	2.11%	0.243%
36	13.139	1721	1726	1734	rVV6	31282	89070	0.94%	0.109%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

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	13.239	1737	1743	1749	rVV	60968	110957	1.17%	0.135%
38	13.498	1784	1787	1793	rVB6	14056	21440	0.23%	0.026%
39	13.651	1808	1813	1814	rBV5	10890	14187	0.15%	0.017%
40	13.698	1815	1821	1828	rVV	1004623	1409785	14.91%	1.722%
41	13.769	1828	1833	1844	rVB3	34212	76444	0.81%	0.093%
42	13.986	1857	1870	1879	rBV	1149763	1982475	20.97%	2.421%
43	14.186	1894	1904	1914	rBV2	66057	196700	2.08%	0.240%
44	14.292	1915	1922	1931	rVV	1296311	2032165	21.49%	2.482%
45	14.504	1951	1958	1959	rBV	1115295	1474427	15.59%	1.801%
46	14.522	1959	1961	1970	rVV	1324316	1610056	17.03%	1.966%
47	14.610	1971	1976	1990	rVV	243021	621251	6.57%	0.759%
48	14.727	1992	1996	2007	rVV3	96260	189895	2.01%	0.232%
49	14.822	2008	2012	2020	rVV5	37550	71932	0.76%	0.088%
50	15.133	2062	2065	2073	rVB6	32019	62293	0.66%	0.076%
51	15.310	2089	2095	2103	rBV	1484104	2244567	23.74%	2.741%
52	15.474	2118	2123	2132	rBV	289652	503926	5.33%	0.615%
53	15.827	2180	2183	2186	rBV4	16512	19766	0.21%	0.024%
54	16.039	2215	2219	2220	rBV4	18697	25116	0.27%	0.031%
55	16.127	2231	2234	2238	rBV2	28418	37471	0.40%	0.046%
56	16.198	2243	2246	2251	rVV6	31676	49958	0.53%	0.061%
57	16.257	2251	2256	2262	rBV	411828	494595	5.23%	0.604%
58	16.504	2293	2298	2313	rBV	391517	786396	8.32%	0.960%
59	16.616	2313	2317	2325	rVV	40825	80228	0.85%	0.098%
60	17.010	2380	2384	2390	rBV3	76483	109826	1.16%	0.134%
61	17.110	2392	2401	2406	rVV	1484632	2464156	26.06%	3.009%
62	17.151	2406	2408	2412	rVB4	162002	170897	1.81%	0.209%
63	17.210	2412	2418	2428	rBV	1603649	2449302	25.90%	2.991%
64	17.292	2429	2432	2434	rBV4	48029	65749	0.70%	0.080%
65	17.316	2434	2436	2441	rVB5	42026	56284	0.60%	0.069%
66	17.480	2460	2464	2466	rBV4	67458	94350	1.00%	0.115%
67	17.521	2467	2471	2474	rBV2	158239	230096	2.43%	0.281%
68	17.751	2505	2510	2511	rBV2	298688	404815	4.28%	0.494%
69	17.774	2511	2514	2531	rVB3	441553	1063739	11.25%	1.299%
70	17.927	2537	2540	2543	rBV3	57927	96914	1.02%	0.118%
71	17.986	2547	2550	2552	rVV3	121622	154963	1.64%	0.189%
72	18.039	2552	2559	2572	rVB	1056497	2313071	24.46%	2.825%
73	18.263	2595	2597	2599	rBV2	51053	50665	0.54%	0.062%
74	18.739	2674	2678	2681	rBV6	121302	215290	2.28%	0.263%
75	18.933	2708	2711	2717	rVB2	188689	273557	2.89%	0.334%
76	19.086	2733	2737	2742	rBV	302680	451924	4.78%	0.552%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

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	19.221	2756	2760	2762	rBV3	205258	304801	3.22%	0.372%
78	19.292	2769	2772	2780	rVB	499493	783580	8.29%	0.957%
79	19.374	2782	2786	2791	rBV	431832	607562	6.43%	0.742%
80	19.539	2811	2814	2816	rBV3	114616	122539	1.30%	0.150%
81	19.592	2819	2823	2841	rVB	2318608	3865583	40.88%	4.720%
82	19.921	2875	2879	2885	rBV2	239368	475868	5.03%	0.581%
83	20.133	2911	2915	2923	rVB2	773477	1509132	15.96%	1.843%
84	20.504	2976	2978	2981	rVB	207597	202693	2.14%	0.248%
85	20.610	2994	2996	3000	rVB2	274744	284236	3.01%	0.347%
86	20.680	3005	3008	3012	rVB2	297569	370882	3.92%	0.453%
87	20.998	3059	3062	3063	rBV	427830	398288	4.21%	0.486%
88	21.021	3063	3066	3070	rVB	682422	1005899	10.64%	1.228%
89	21.198	3091	3096	3106	rVB	4035049	6589258	69.69%	8.047%
90	21.551	3151	3156	3165	rBV2	1599762	3258653	34.46%	3.979%
91	22.109	3249	3251	3253	rBV	231178	163992	1.73%	0.200%
92	22.174	3259	3262	3263	rBV2	367602	416035	4.40%	0.508%
93	22.427	3296	3305	3312	rBV	4288992	9455350	100.00%	11.546%
94	23.480	3481	3484	3491	rVB	468774	692705	7.33%	0.846%
95	23.939	3558	3562	3570	rVB	1330280	3170720	33.53%	3.872%
96	24.021	3571	3576	3591	rVB2	1331517	3675775	38.88%	4.489%
97	24.645	3675	3682	3691	rVB	1296346	3602049	38.10%	4.399%
98	24.880	3716	3722	3729	rBV	1182807	3074033	32.51%	3.754%
99	26.262	3953	3957	3971	rVB2	542954	1644303	17.39%	2.008%

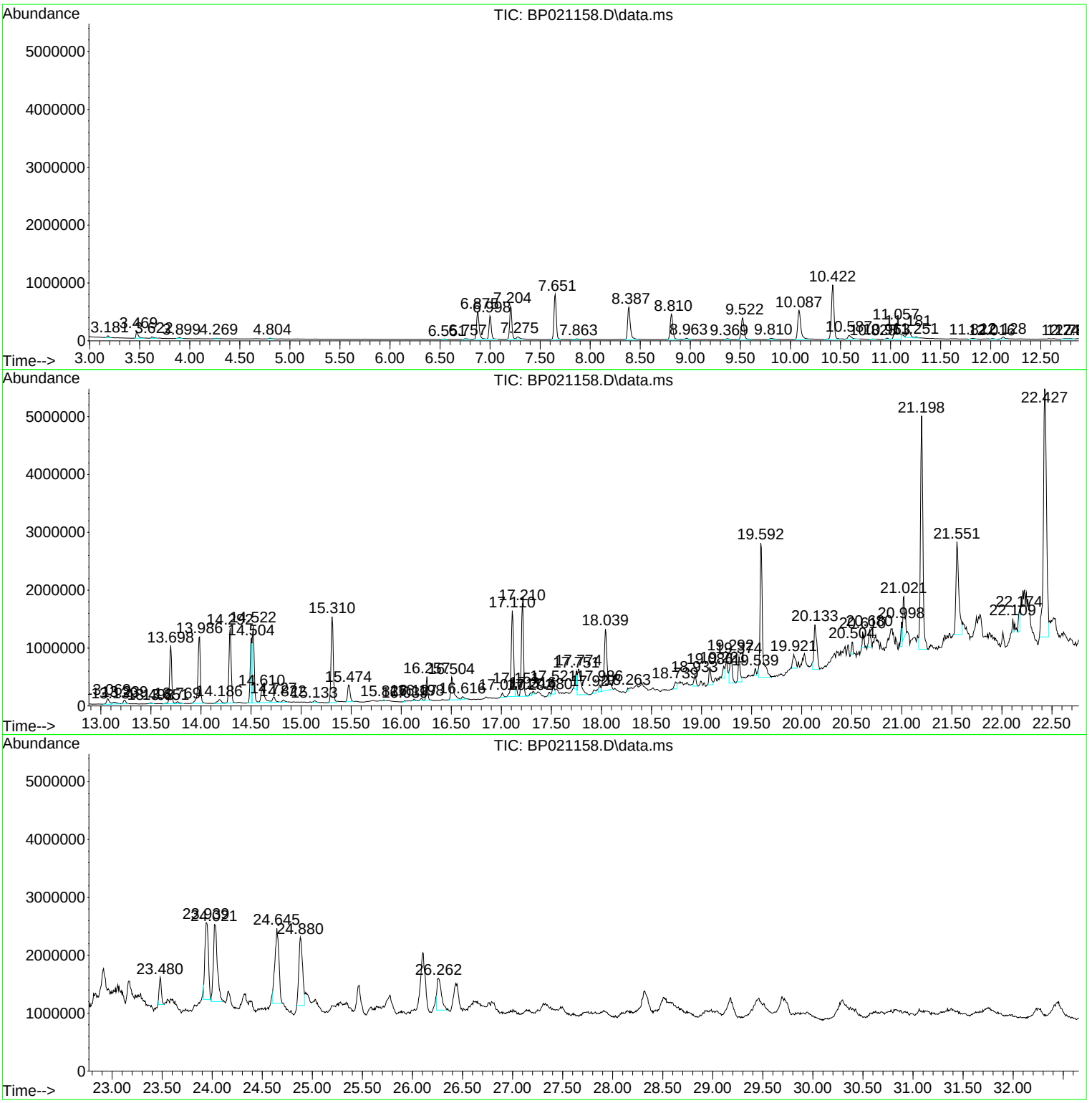
Sum of corrected areas: 81889582

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

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\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

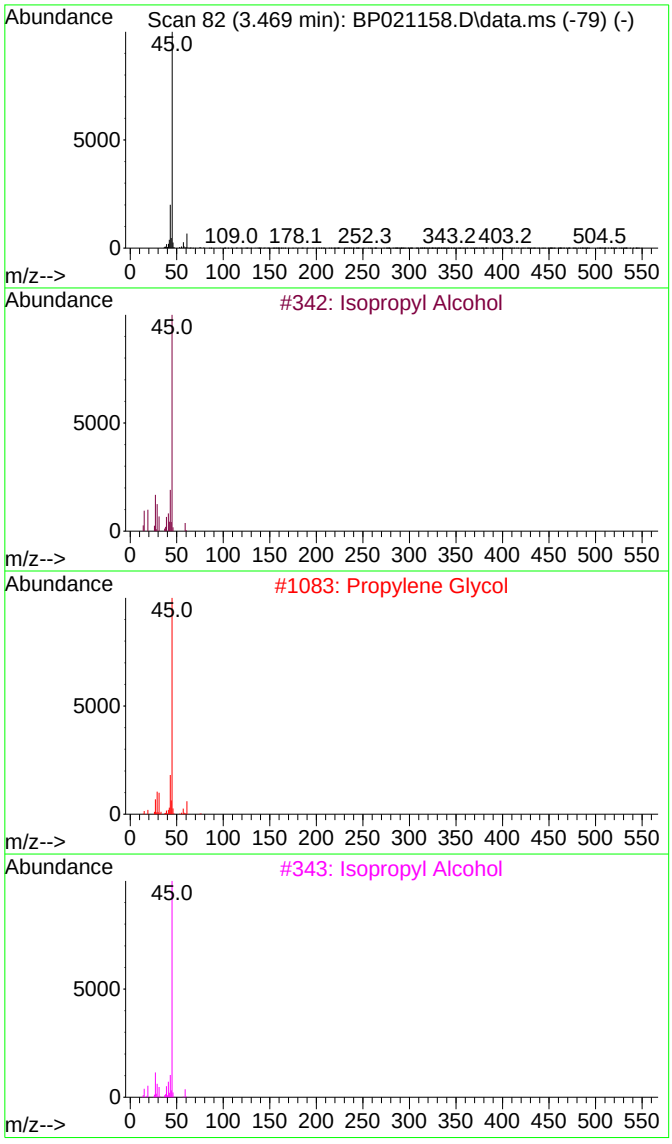
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

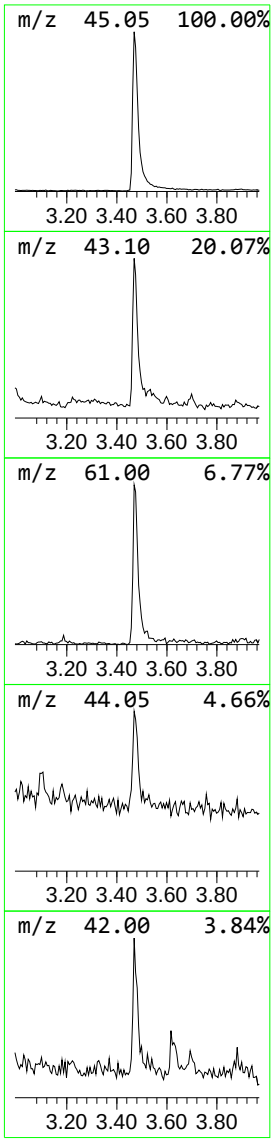
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Isopropyl Alcohol Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.469	2.93 ng/ul	176947	1,4-Dichlorobenzene-d4	7.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
2	Propylene Glycol	76	C3H8O2	000057-55-6	9
3	Isopropyl Alcohol	60	C3H8O	000067-63-0	9
4	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
5	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

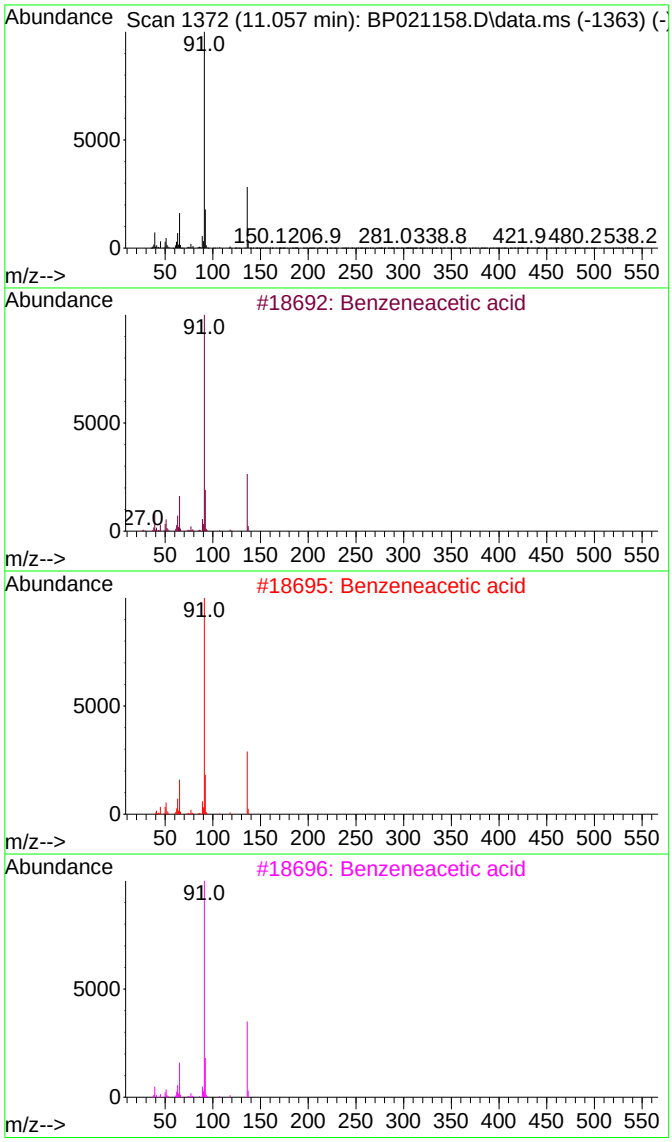
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

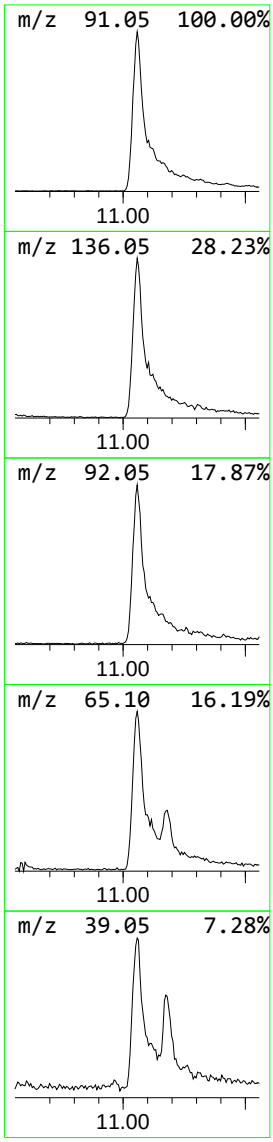
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzeneacetic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.057	9.86 ng/ul	834762	Naphthalene-d8	10.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetic acid	136	C8H8O2	000103-82-2	95
2	Benzeneacetic acid	136	C8H8O2	000103-82-2	94
3	Benzeneacetic acid	136	C8H8O2	000103-82-2	94
4	Benzeneacetic acid	136	C8H8O2	000103-82-2	90
5	Propanedioic acid, phenyl-	180	C9H8O4	002613-89-0	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

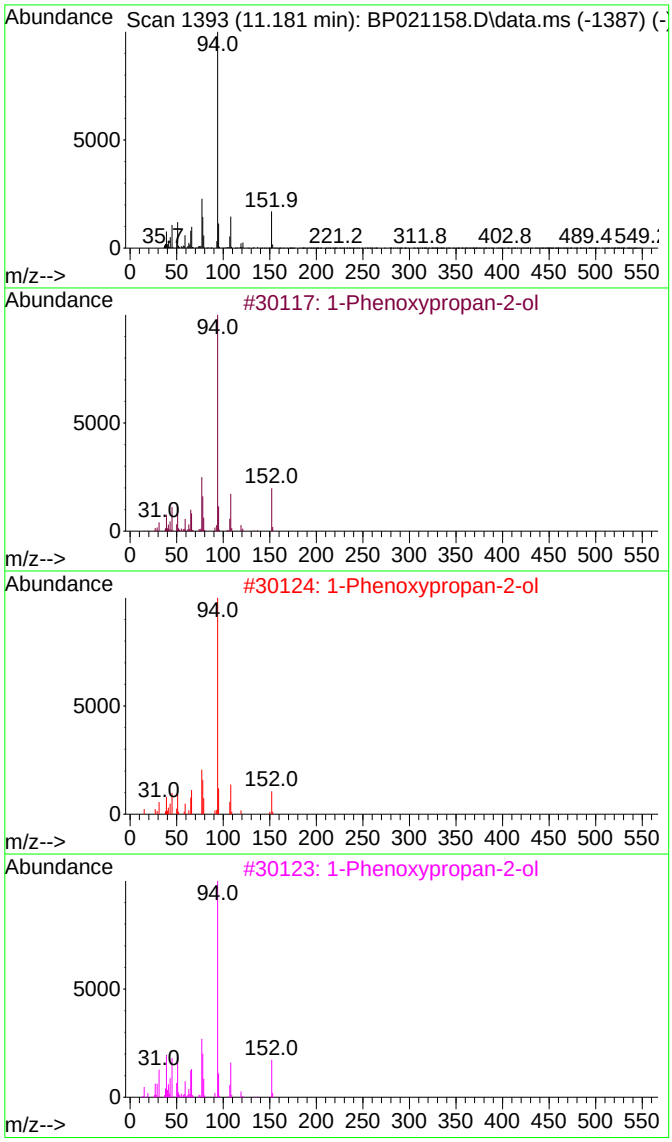
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

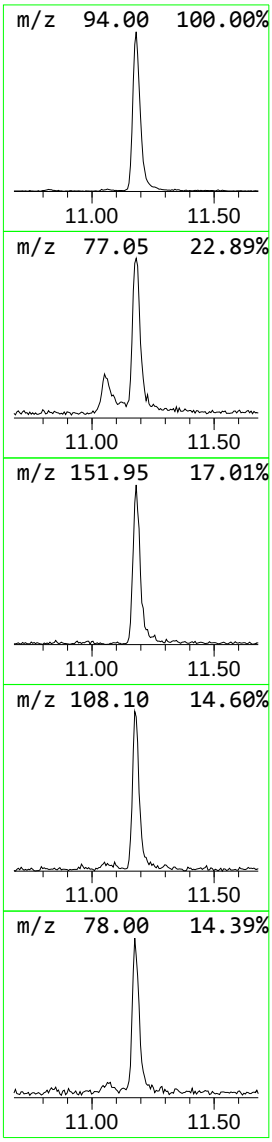
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.181	3.82 ng/ul	323050	Naphthalene-d8	10.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
5	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

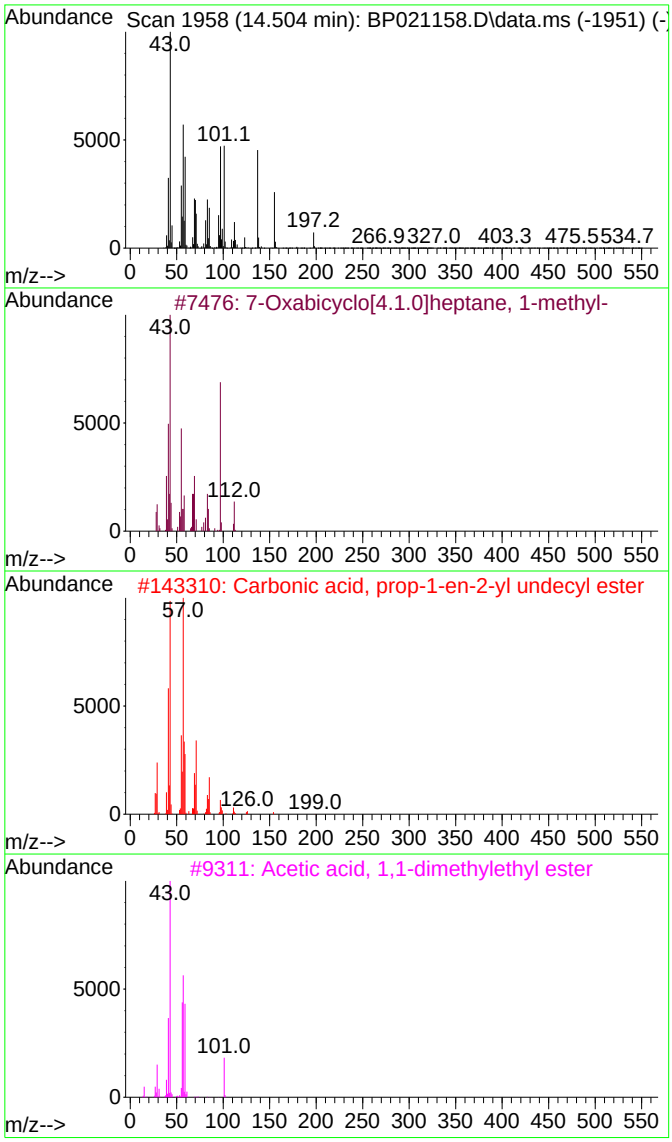
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

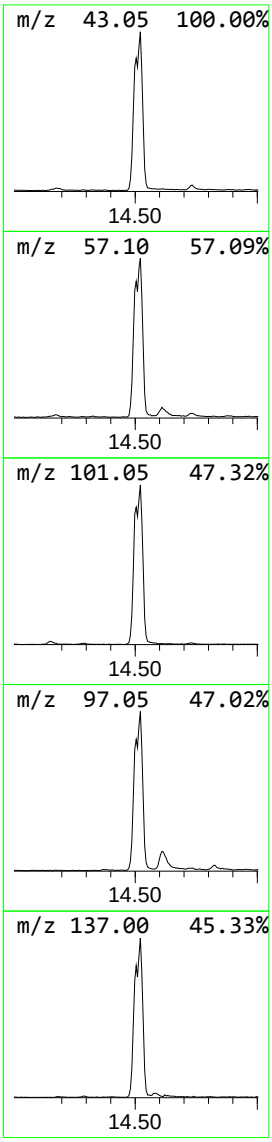
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.504	14.51 ng/ul	1474430	Acenaphthene-d10	14.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Oxabicyclo[4.1.0]heptane, 1-me...	112	C7H12O	001713-33-3	10
2	Carbonic acid, prop-1-en-2-yl un...	256	C15H28O3	1000382-90-6	10
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	10
4	trans-5-Isopropyl-6,7-epoxy-8-hy...	228	C13H24O3	058002-07-6	10
5	Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	10





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

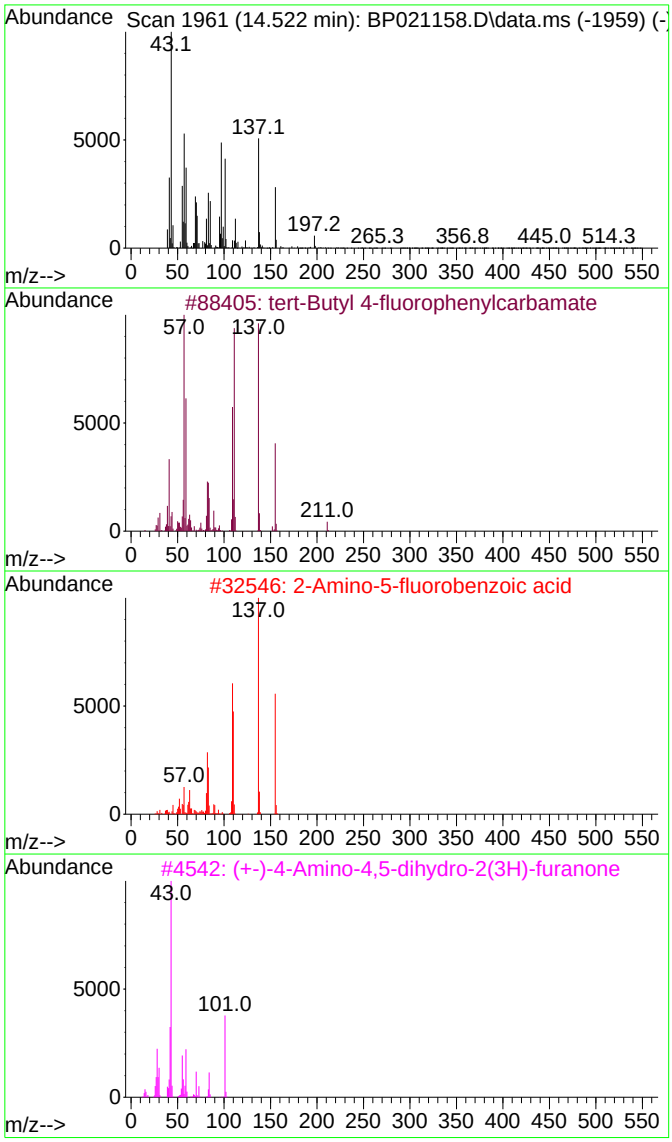
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

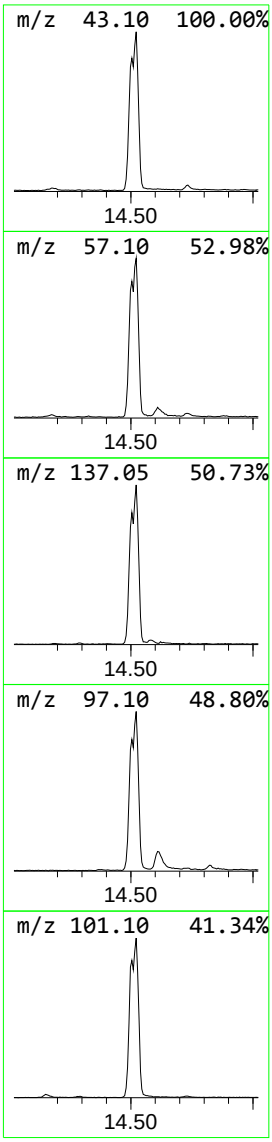
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.522	15.85 ng/ul	1610060	Acenaphthene-d10	14.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	tert-Butyl 4-fluorophenylcarbamate	211	C11H14FNO2	060144-53-8	27
2	2-Amino-5-fluorobenzoic acid	155	C7H6FNO2	000446-08-2	18
3	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	14
4	Decane, 1-bromo-	220	C10H21Br	000112-29-8	10
5	2,2,3,4-Tetramethylhex-5-en-3-ol	156	C10H20O	1010191-06-9	10





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Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

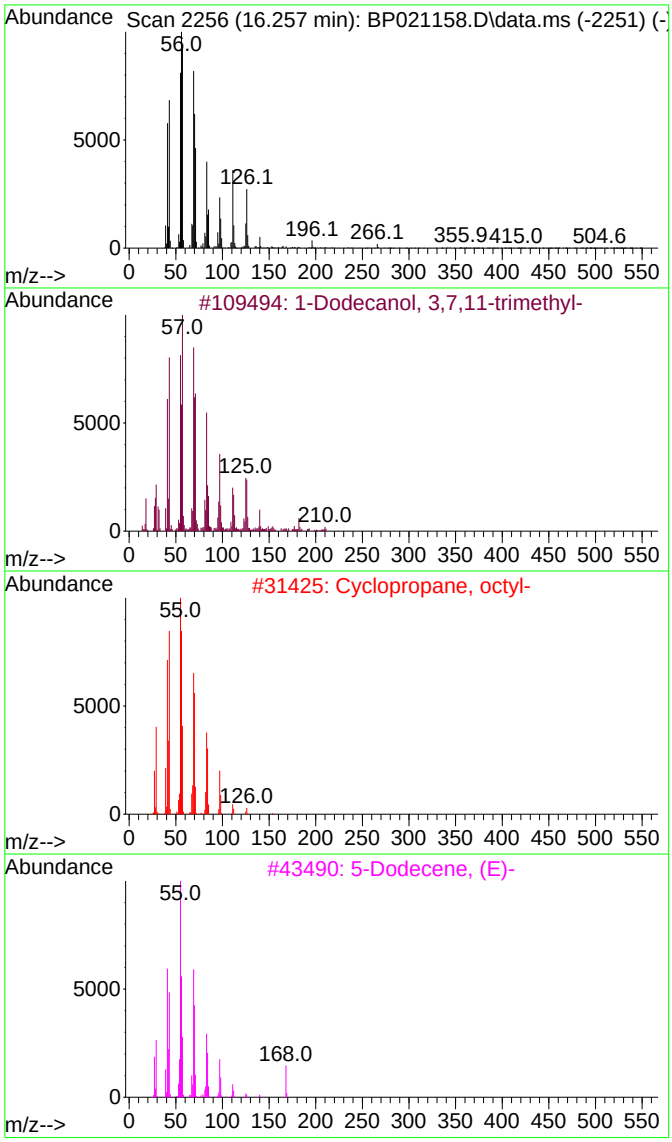
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

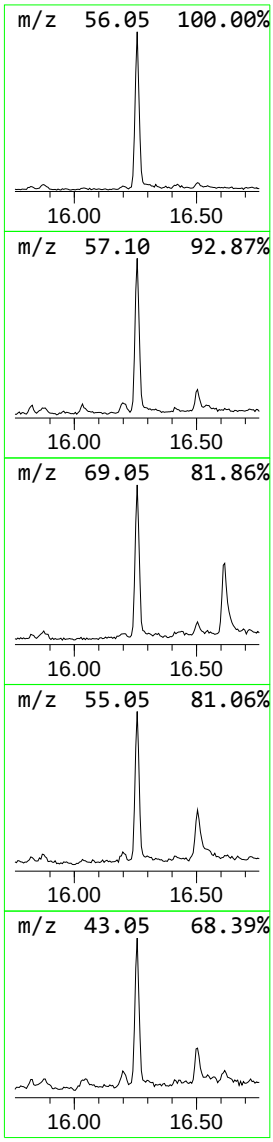
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1-Dodecanol, 3,7,11-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.257	4.01 ng/ul	494595	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Dodecanol, 3,7,11-trimethyl-	228	C15H32O	006750-34-1	72
2	Cyclopropane, octyl-	154	C11H22	001472-09-9	64
3	5-Dodecene, (E)-	168	C12H24	007206-16-8	58
4	1-Undecene, 8-methyl-	168	C12H24	074630-40-3	58
5	5-Dodecene, (E)-	168	C12H24	007206-16-8	58





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Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
A4CG7

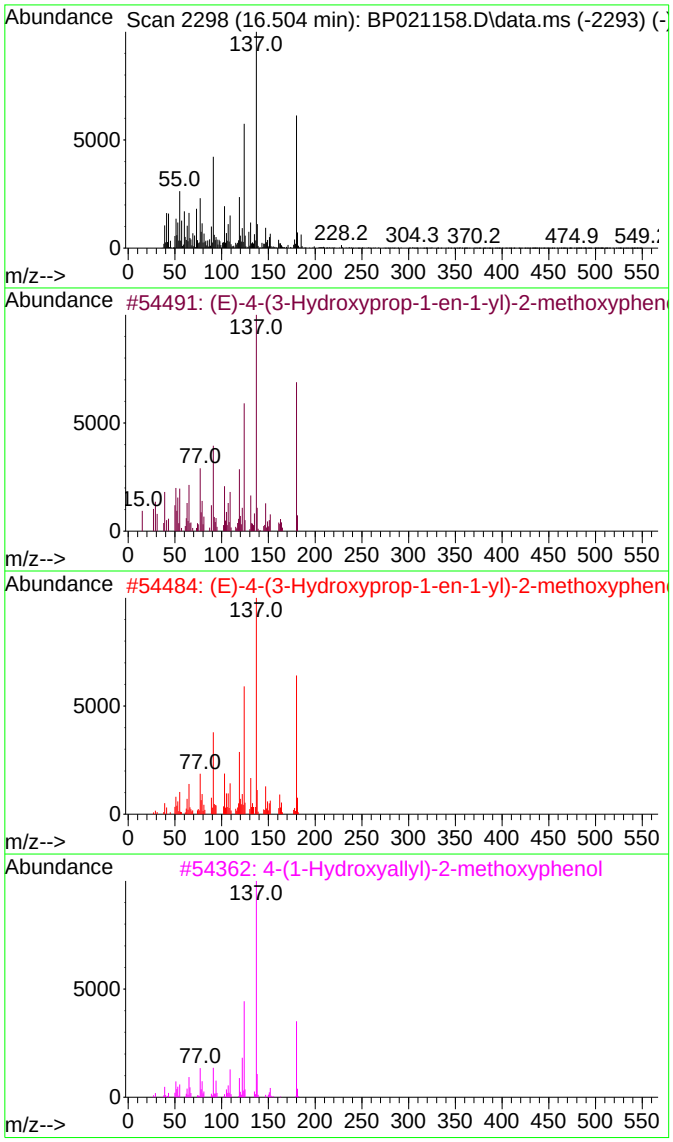
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

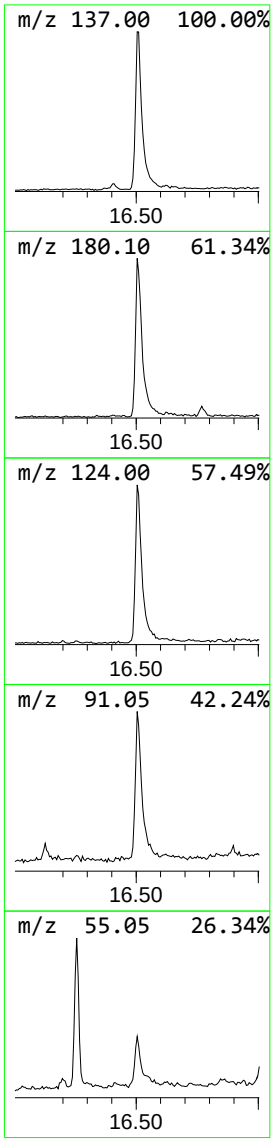
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (E)-4-(3-Hydroxyprop-1-en-1... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.504	6.38 ng/ul	786396	Phenanthrene-d10	17.110

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	Guaiacol, 4-butyl-	180	C11H16O2	059832-96-1	46
5	5-(Acetylaminomethyl)-4-amino-2-...	180	C8H12N4O	023676-63-3	43





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Misc :
ALS Vial : 18 Sample Multiplier: 1

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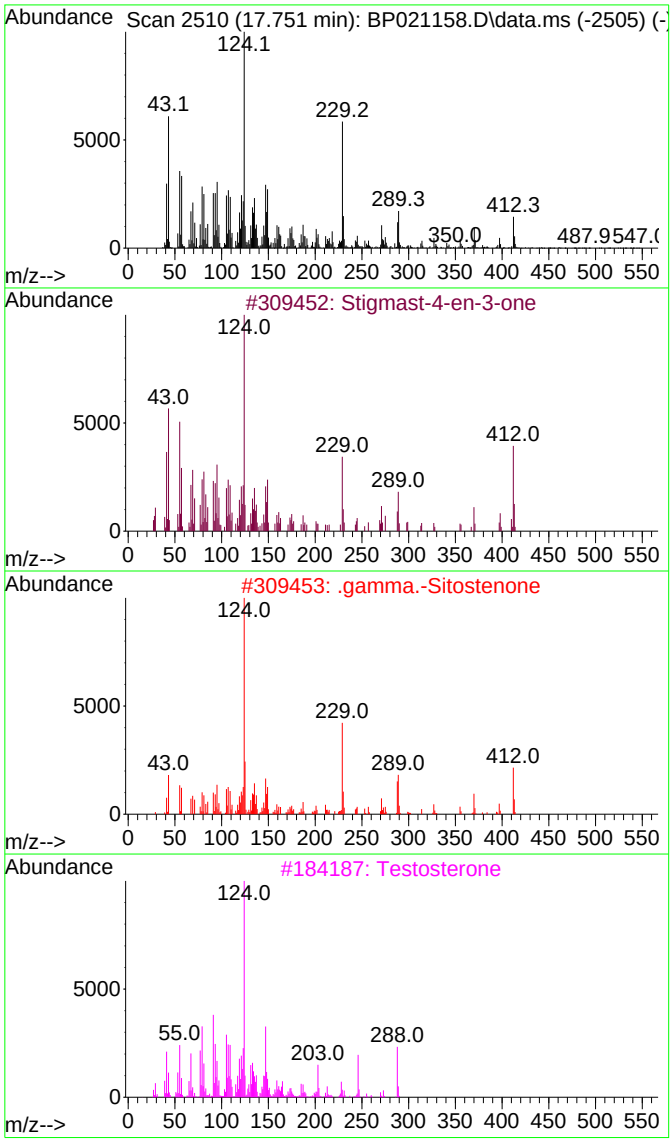
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

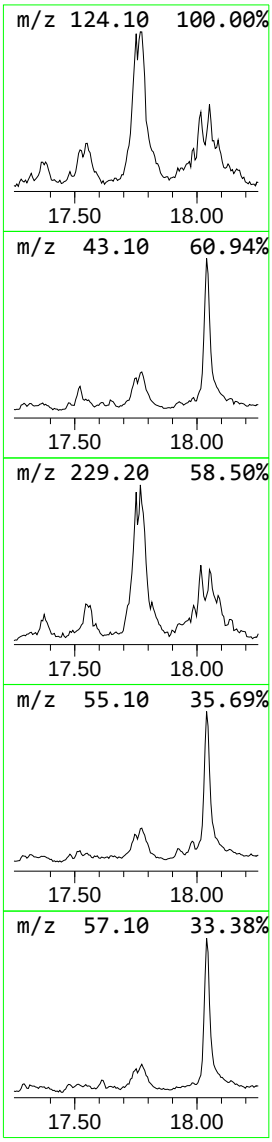
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Stigmast-4-en-3-one Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.751	3.29 ng/ul	404815	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Stigmast-4-en-3-one	412	C29H48O	001058-61-3	95
2	.gamma.-Sitostenone	412	C29H48O	084924-96-9	90
3	Testosterone	288	C19H28O2	000058-22-0	86
4	Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000604-39-7	60
5	Testosterone	288	C19H28O2	000058-22-0	46





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Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

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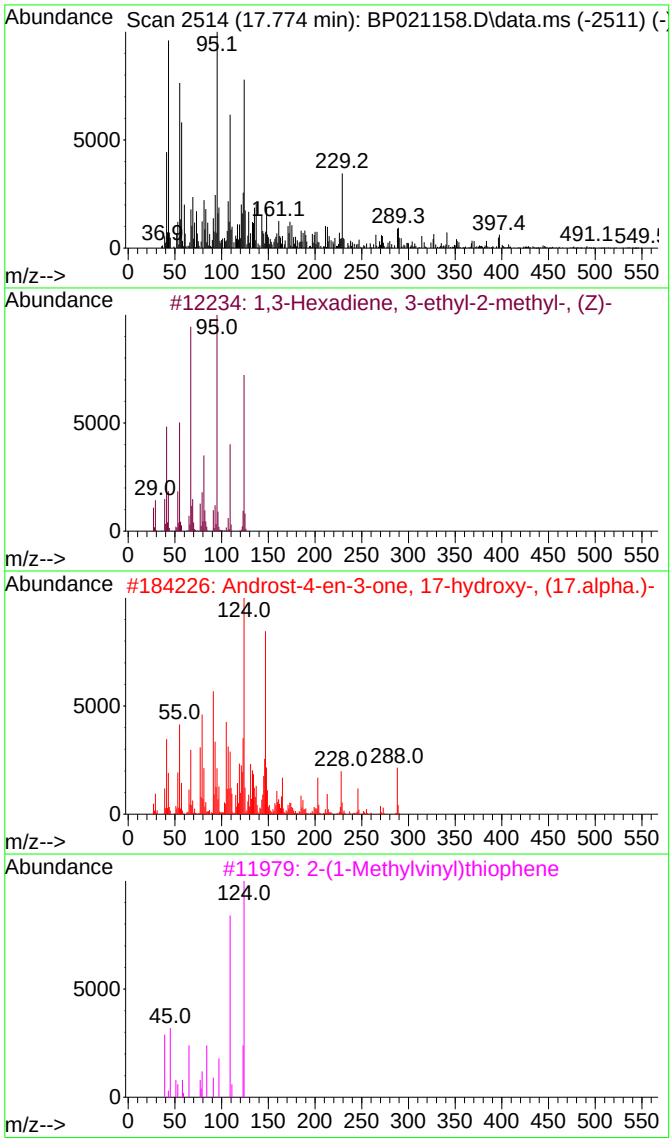
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

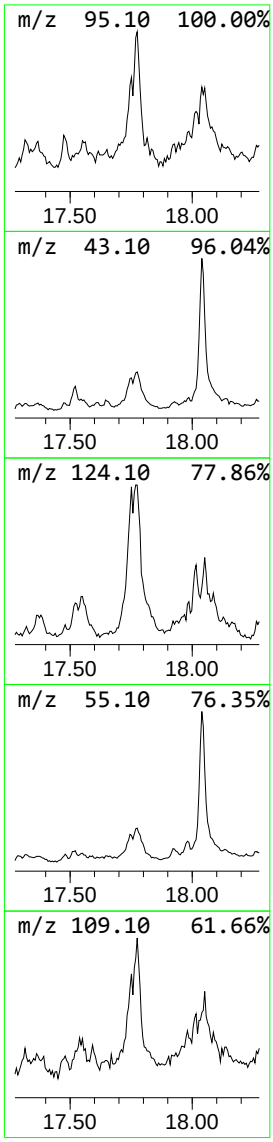
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1,3-Hexadiene, 3-ethyl-2-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.774	8.63 ng/ul	1063740	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Hexadiene, 3-ethyl-2-methyl-...	124	C9H16	074752-97-9	50
2	Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000481-30-1	43
3	2-(1-Methylvinyl)thiophene	124	C7H8S	030616-73-0	30
4	4-Methylimidazole-5-butyric acid	168	C8H12N2O2	1000129-14-7	27
5	1,2-Benzenediol, 3-methyl-	124	C7H8O2	000488-17-5	22





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Acq On : 25 Jul 2024 22:26
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Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

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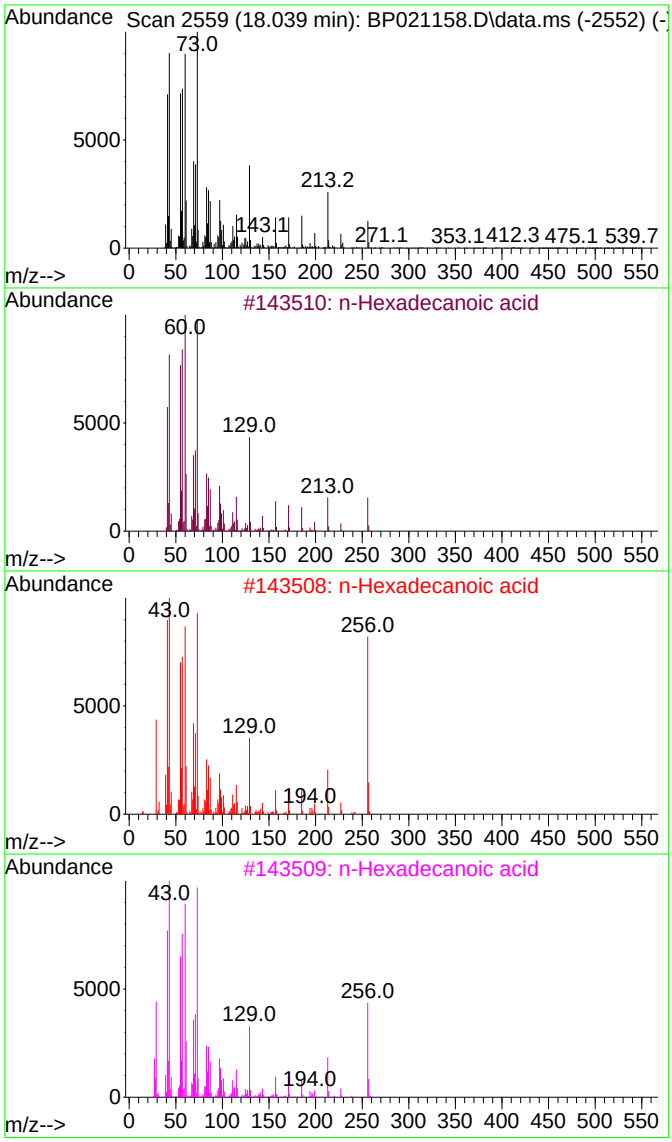
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

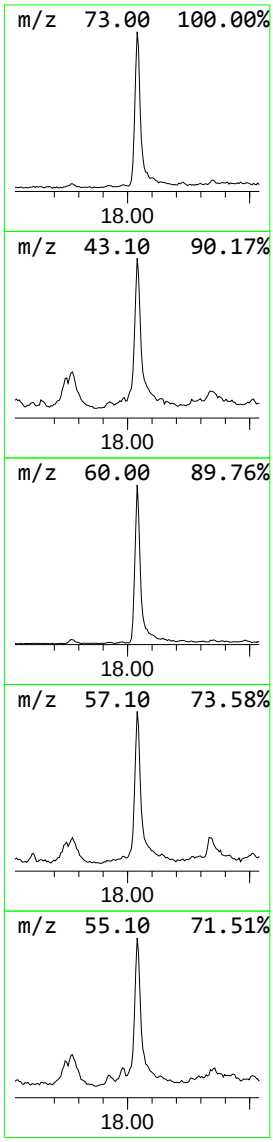
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.039	18.77 ng/ul	2313070	Phenanthrene-d10	17.110

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	98
5	Tridecanoic acid	214	C13H26O2	000638-53-9	95





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Acq On : 25 Jul 2024 22:26
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Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

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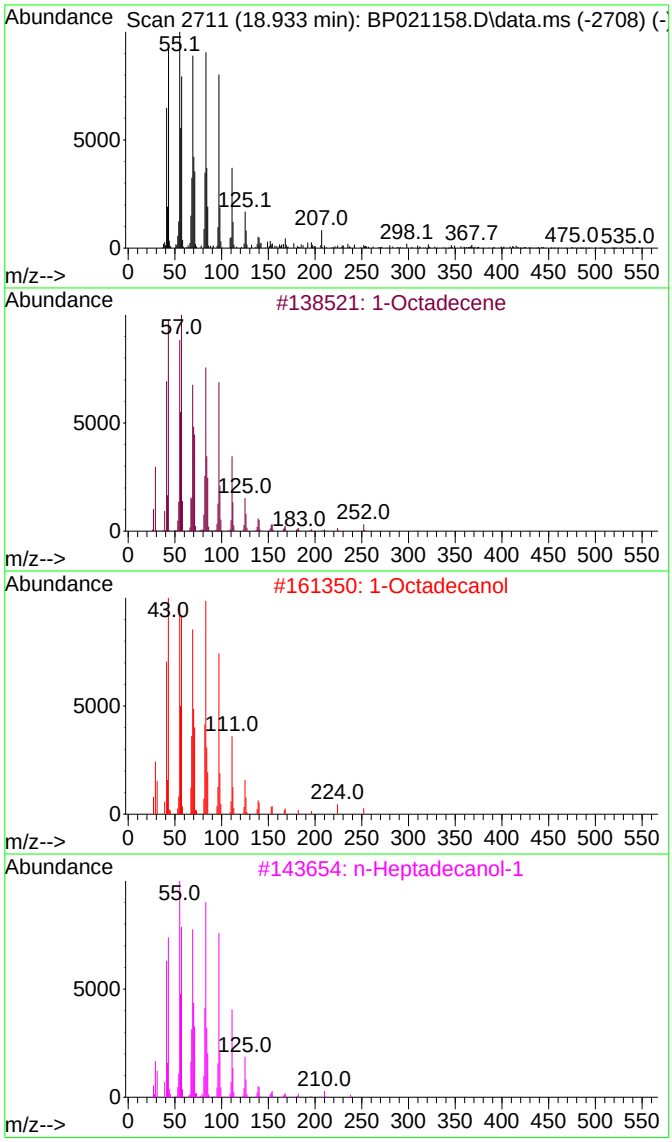
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

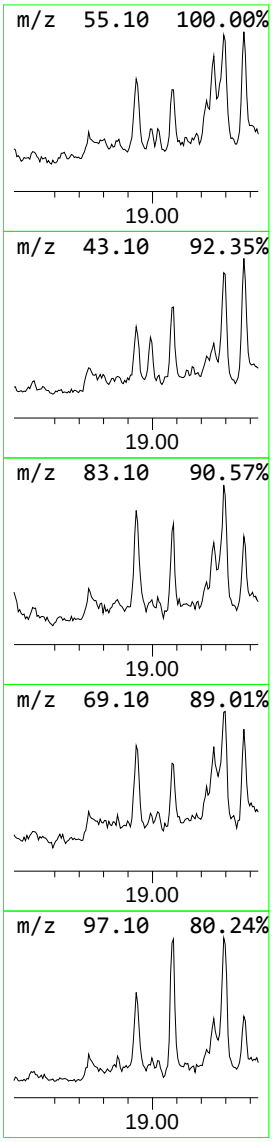
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.933	2.22 ng/ul	273557	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	95
2	1-Octadecanol	270	C18H38O	000112-92-5	94
3	n-Heptadecanol-1	256	C17H36O	001454-85-9	94
4	Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	93





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

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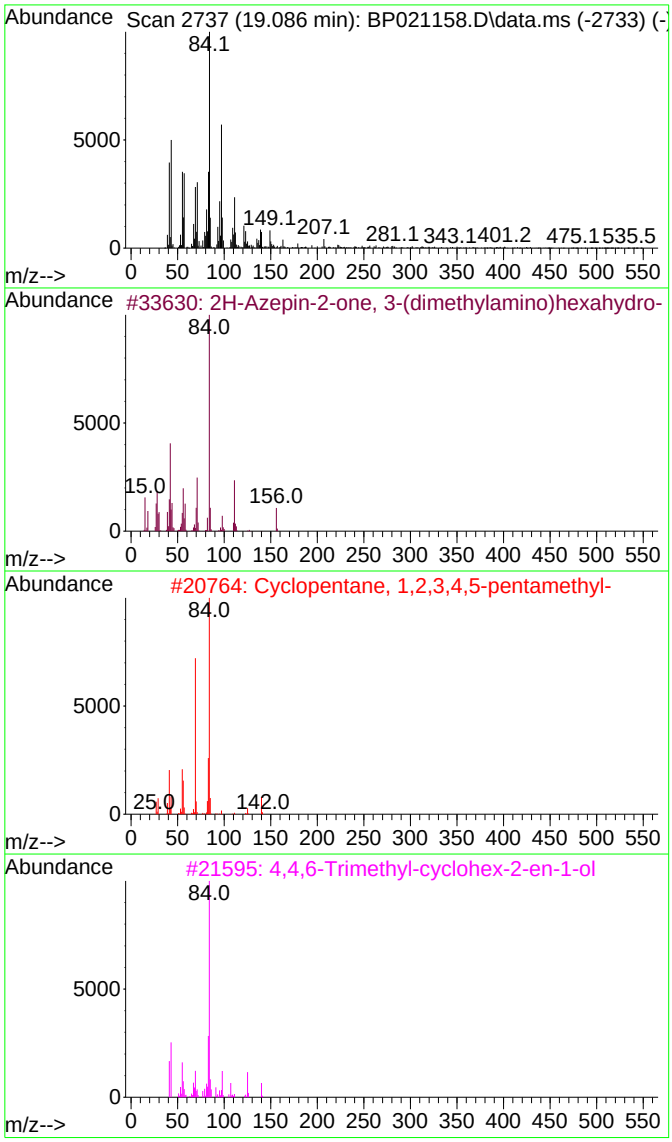
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

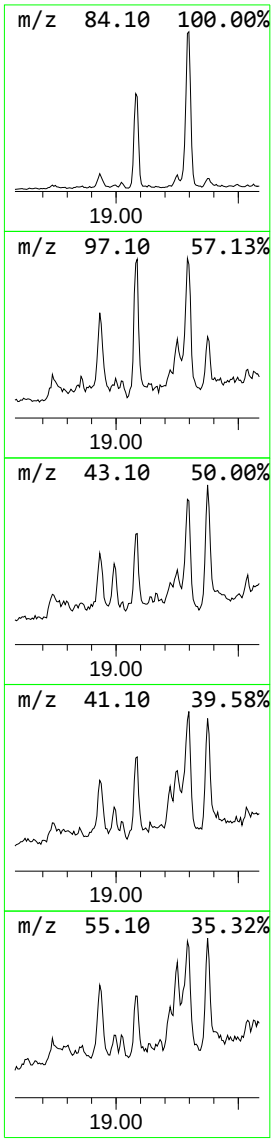
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2H-Azepin-2-one, 3-(dimethy... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.086	3.67 ng/ul	451924	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Azepin-2-one, 3-(dimethylamin...	156	C8H16N2O	070394-23-9	55
2	Cyclopentane, 1,2,3,4,5-pentamet...	140	C10H20	1000152-79-7	38
3	4,4,6-Trimethyl-cyclohex-2-en-1-ol	140	C9H16O	1000144-64-7	38
4	(S)-5-Hydroxymethyl-2[5H]-furanone	114	C5H6O3	078508-96-0	30
5	2-Methoxy-4,4,5,5-tetramethyl-2-...	157	C8H15NO2	049680-47-9	27





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Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

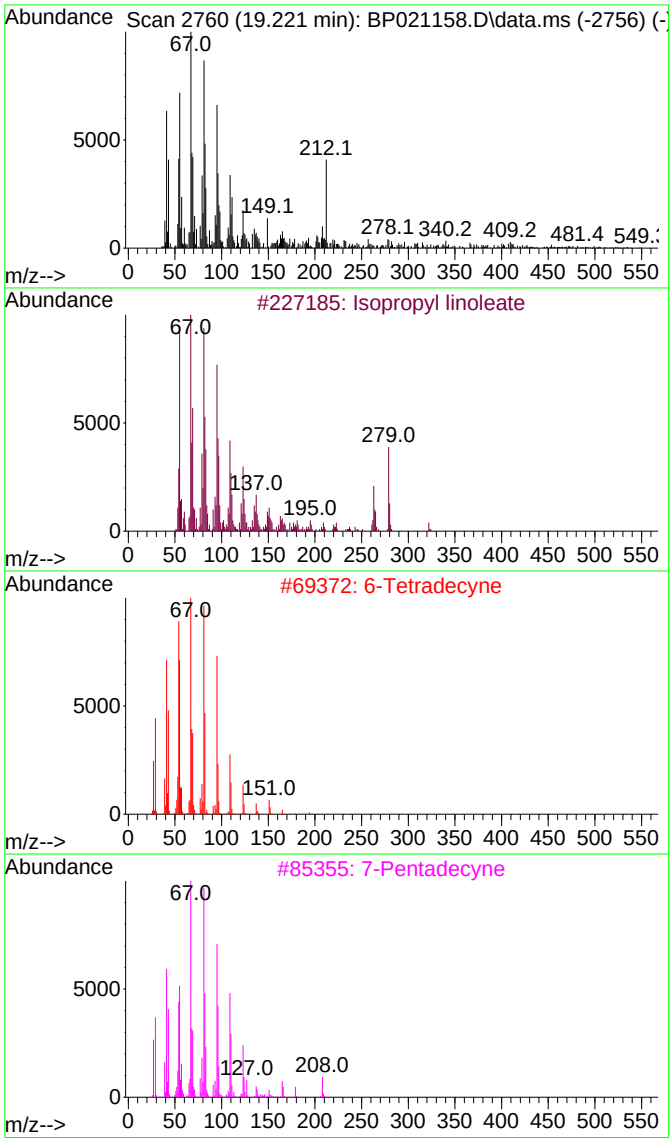
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

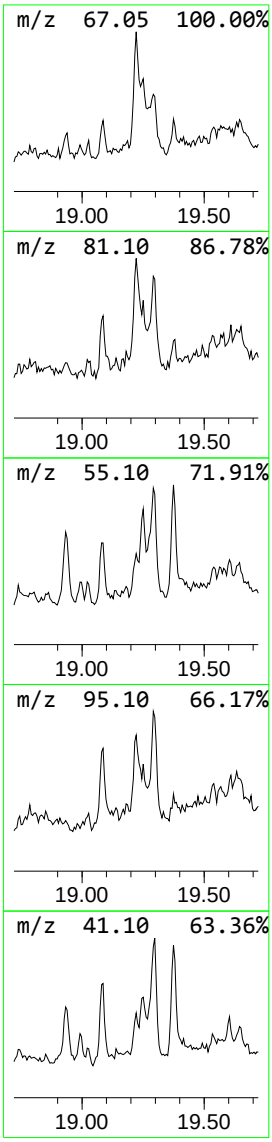
****

Peak Number 13 Isopropyl linoleate Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.221	2.47 ng/ul	304801	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl linoleate	322	C21H38O2	022882-95-7	97
2	6-Tetradecyne	194	C14H26	003730-08-3	89
3	7-Pentadecyne	208	C15H28	022089-89-0	68
4	Linoelaidic acid	280	C18H32O2	000506-21-8	68
5	Linoleyl methyl ketone	278	C19H34O	029204-24-8	64





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Misc :
ALS Vial : 18 Sample Multiplier: 1

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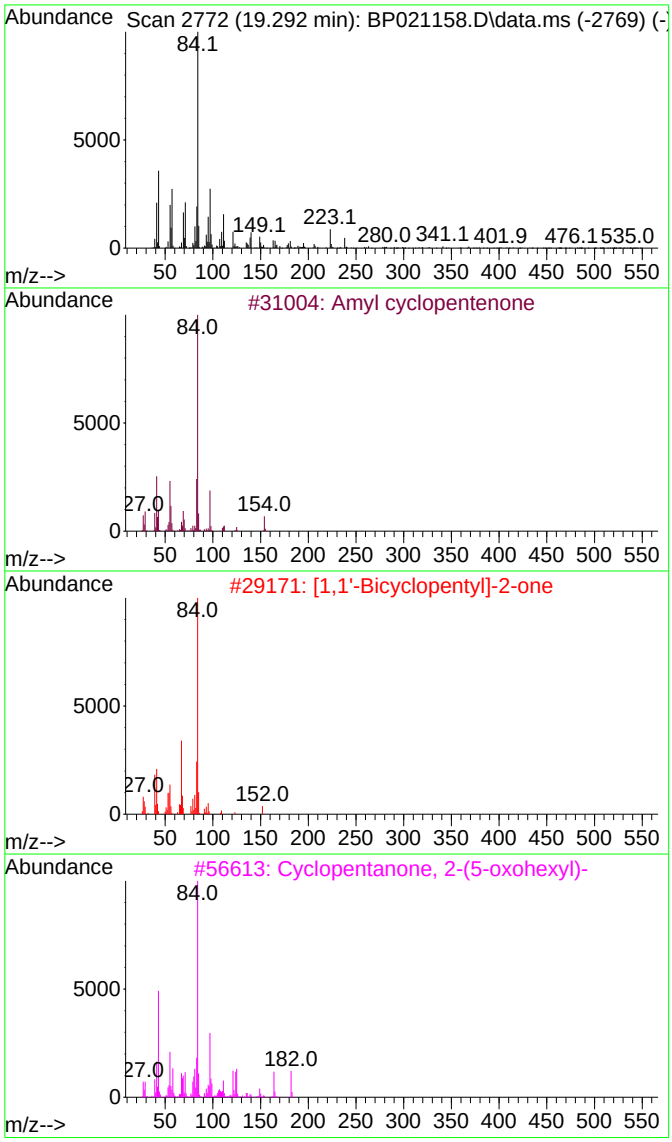
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

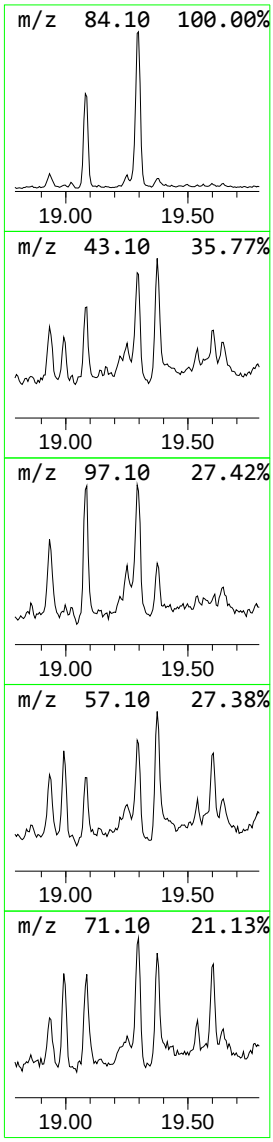
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Amyl cyclopentenone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.292	6.36 ng/ul	783580	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Amyl cyclopentenone	154	C10H18O	004819-67-4	50
2	[1,1'-Bicyclopentyl]-2-one	152	C10H16O	004884-24-6	47
3	Cyclopentanone, 2-(5-oxohexyl)-	182	C11H18O2	015674-93-8	45
4	Cyclohexylamine, N-ethyl-	127	C8H17N	005459-93-8	43
5	2-Cyclohexylpiperidine	167	C11H21N	056528-77-9	43





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ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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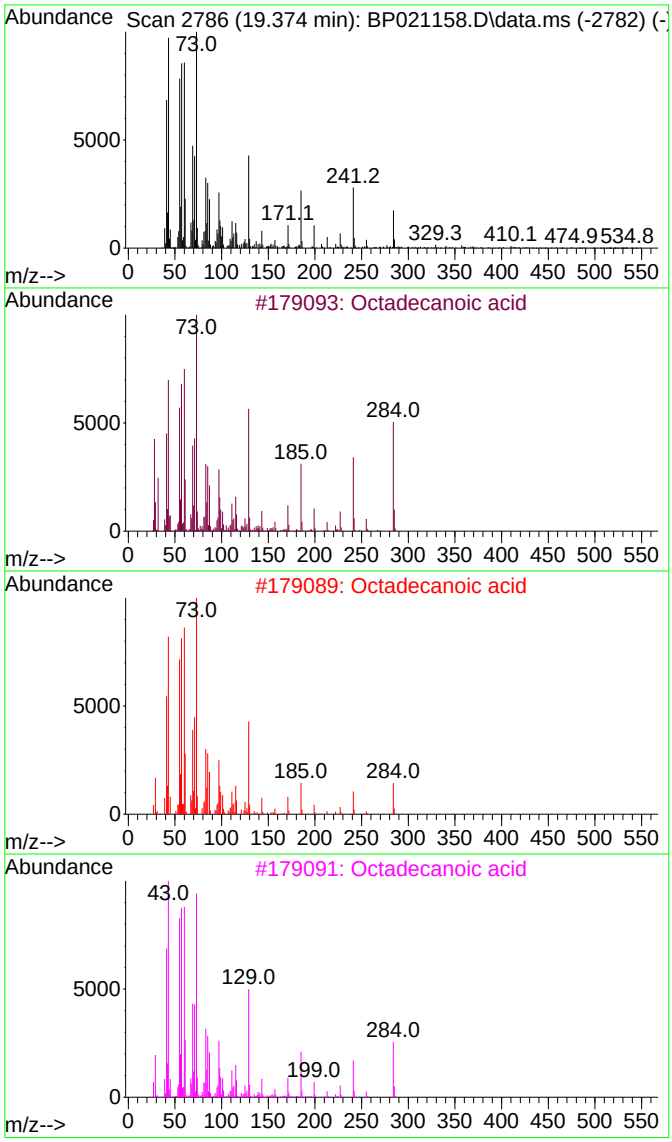
Quant Title : SVOA CALIBRATION

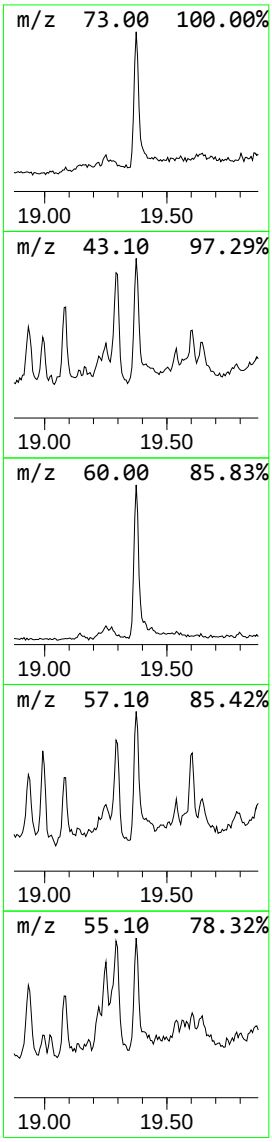
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Octadecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.374	3.73 ng/ul	607562	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	95
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94





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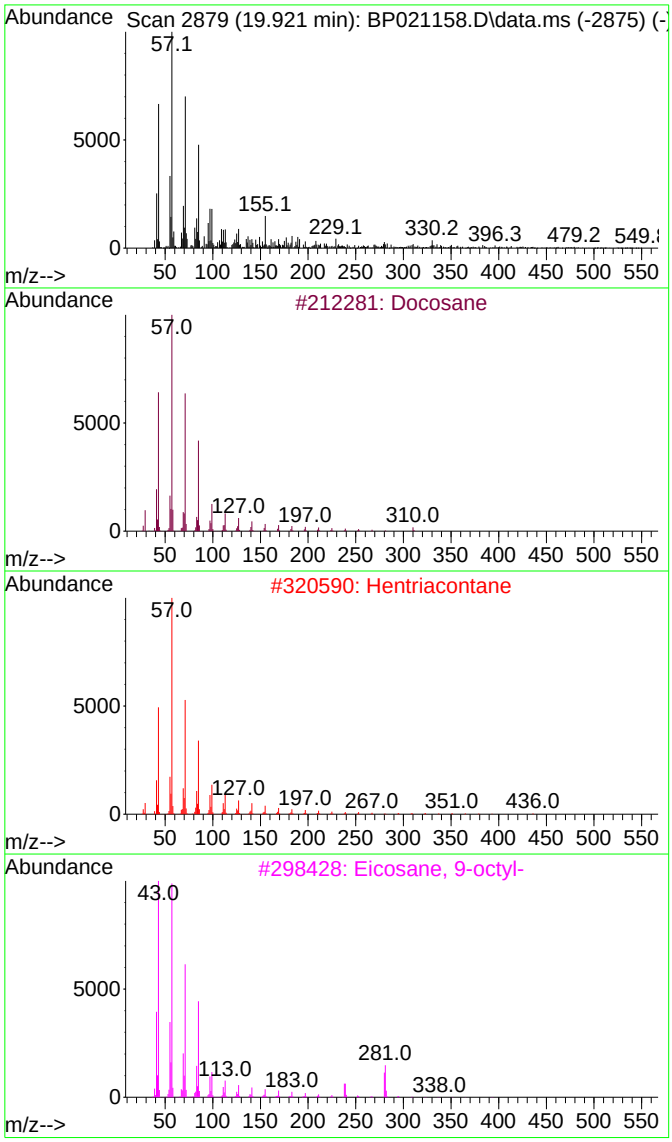
Quant Title : SVOA CALIBRATION

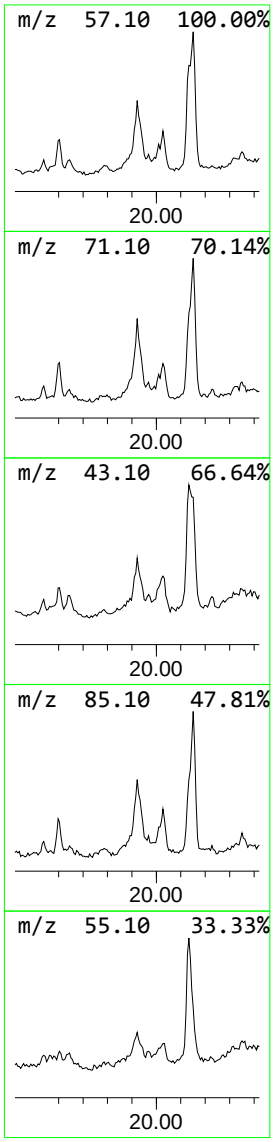
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.921	2.92 ng/ul	475868	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane	310	C22H46	000629-97-0	93
2	Hentriacontane	437	C31H64	000630-04-6	91
3	Eicosane, 9-octyl-	394	C28H58	013475-77-9	90
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	89
5	Nonacosane	408	C29H60	000630-03-5	81





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Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
A4CG7

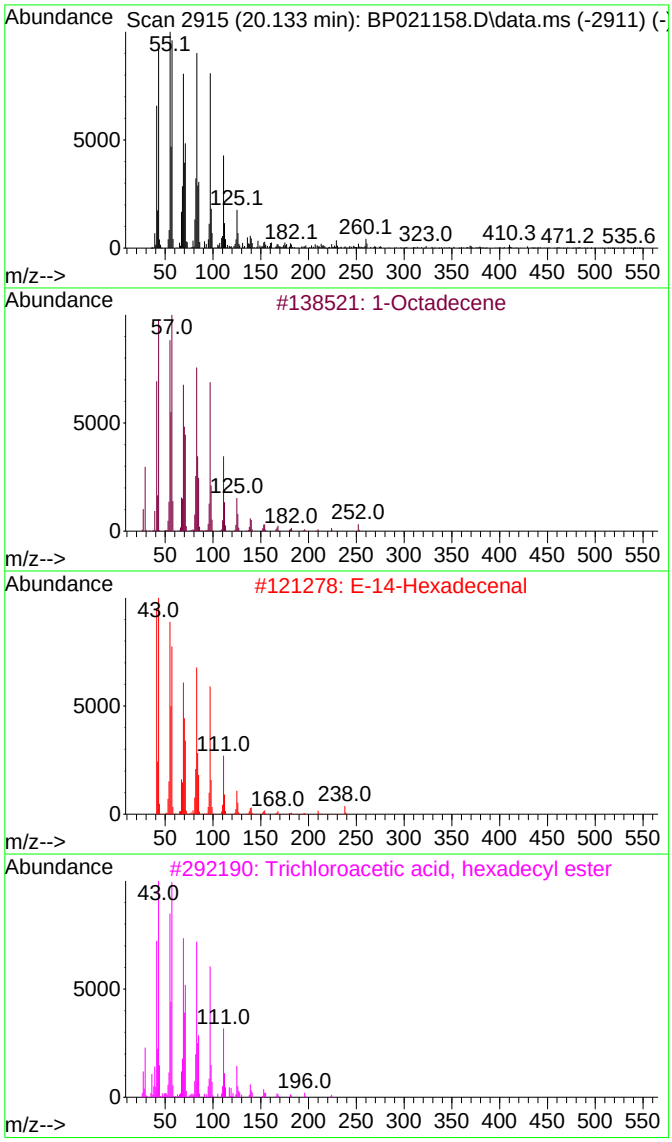
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

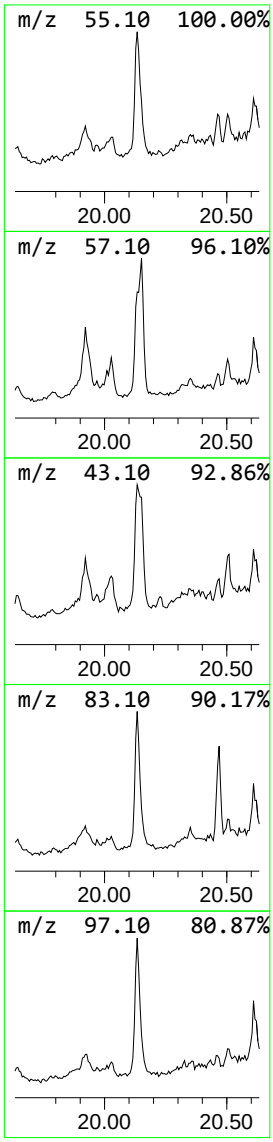
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.133	9.26 ng/ul	1509130	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	98
2	E-14-Hexadecenal	238	C16H30O	330207-53-9	95
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91





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Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

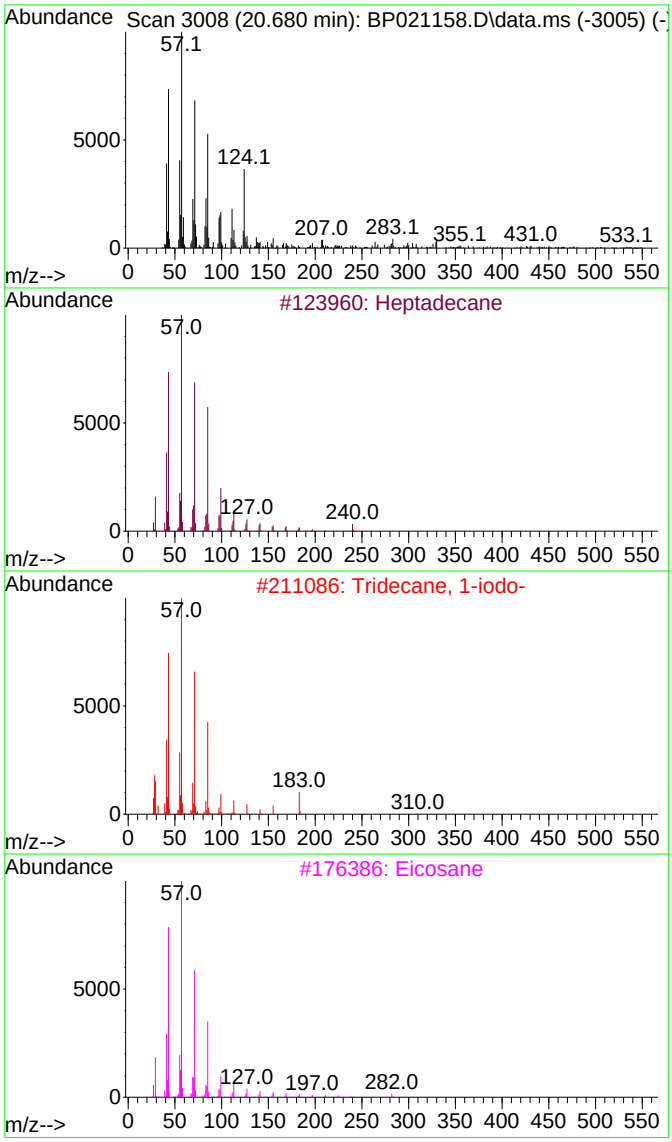
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

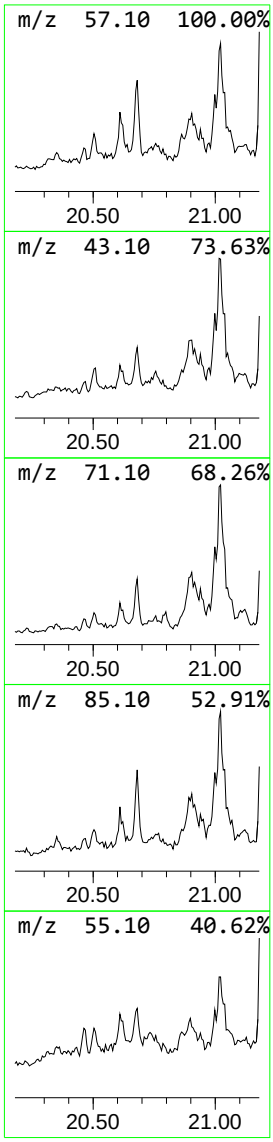
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.680	2.28 ng/ul	370882	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	92
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
3	Eicosane	282	C20H42	000112-95-8	86
4	Dotriacontane, 2-methyl-	465	C33H68	001720-11-2	83
5	Eicosane	282	C20H42	000112-95-8	64





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

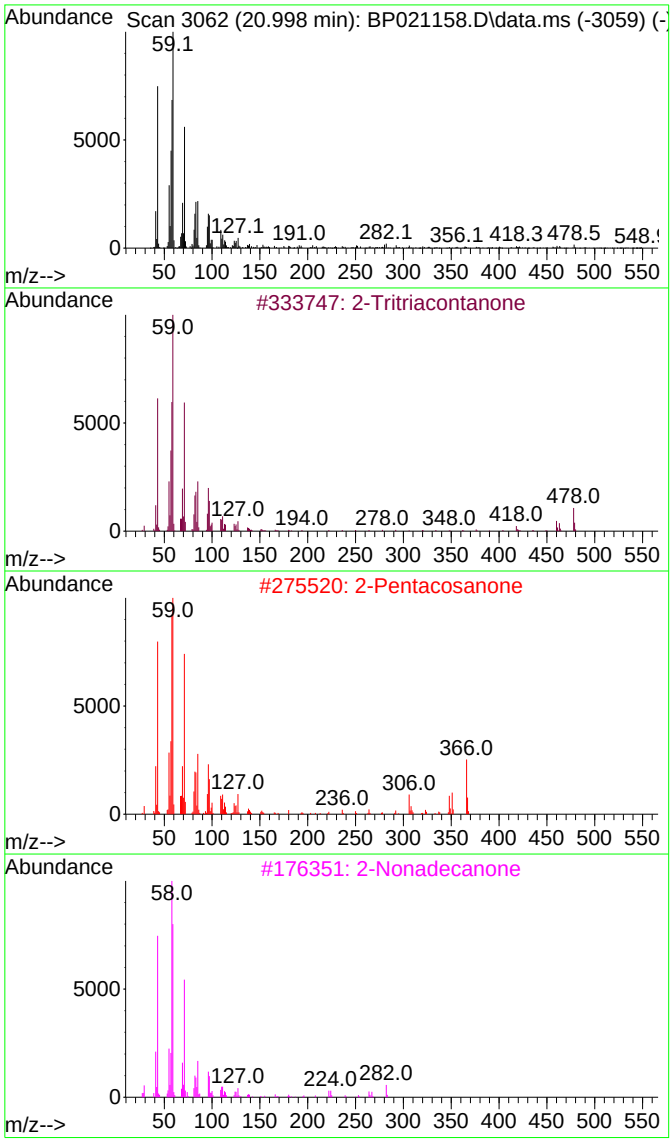
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

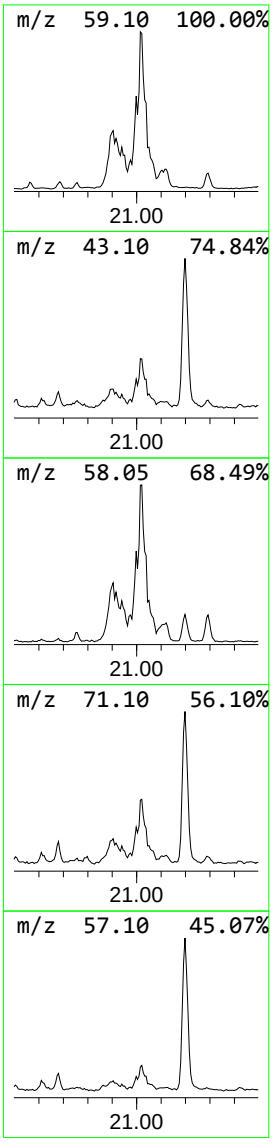
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 2-Tritriacontanone Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.998	2.44 ng/ul	398288	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Tritriacontanone	479	C33H66O	075207-55-5	98
2	2-Pentacosanone	366	C25H50O	075207-54-4	93
3	2-Nonadecanone	282	C19H38O	000629-66-3	76
4	Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	52
5	Carbonic acid, 2-methoxyethyl ne...	190	C9H18O4	1000331-42-7	50





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

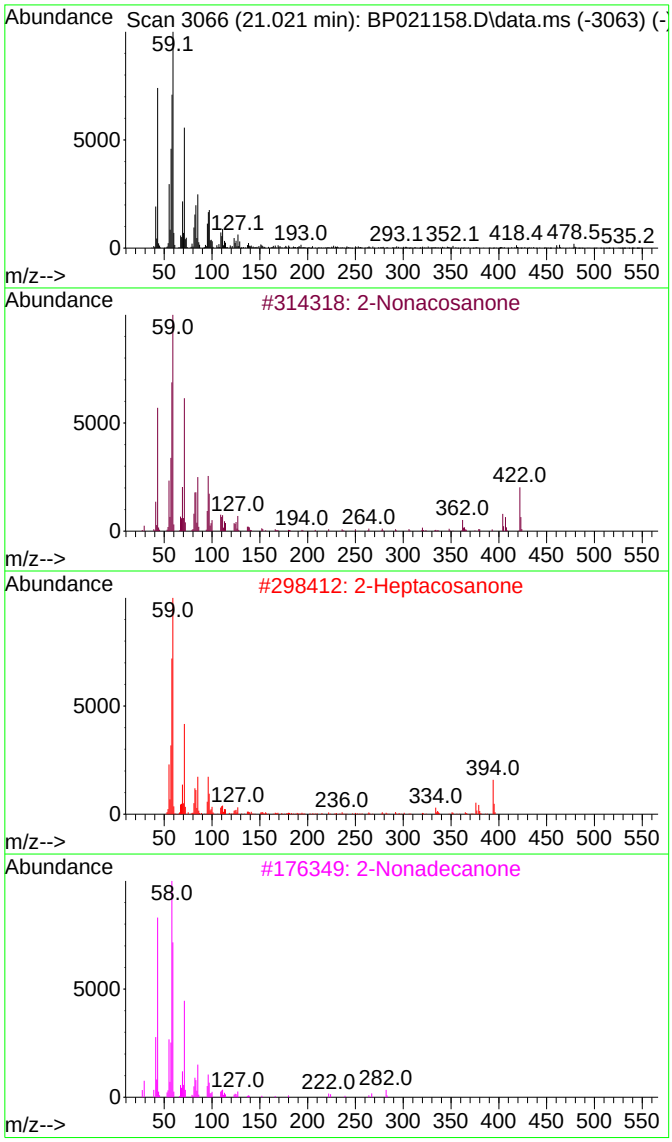
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

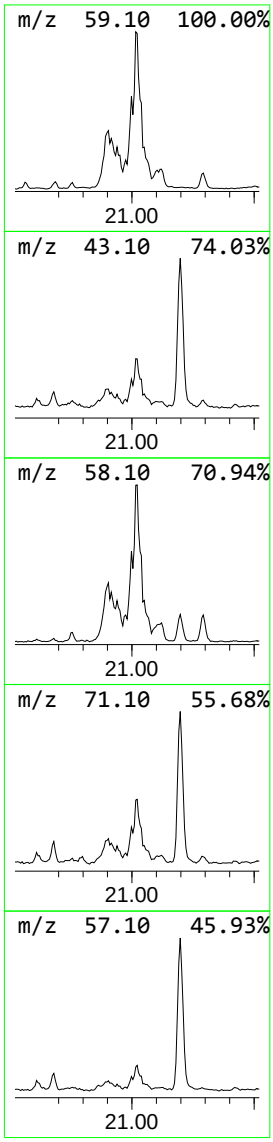
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 2-Nonacosanone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.021	6.17 ng/ul	1005900	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Nonacosanone	422	C29H58O	017600-99-6	91
2	2-Heptacosanone	394	C27H54O	007796-19-2	83
3	2-Nonadecanone	282	C19H38O	000629-66-3	64
4	Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	53
5	2-Pentadecanone	226	C15H30O	002345-28-0	43





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

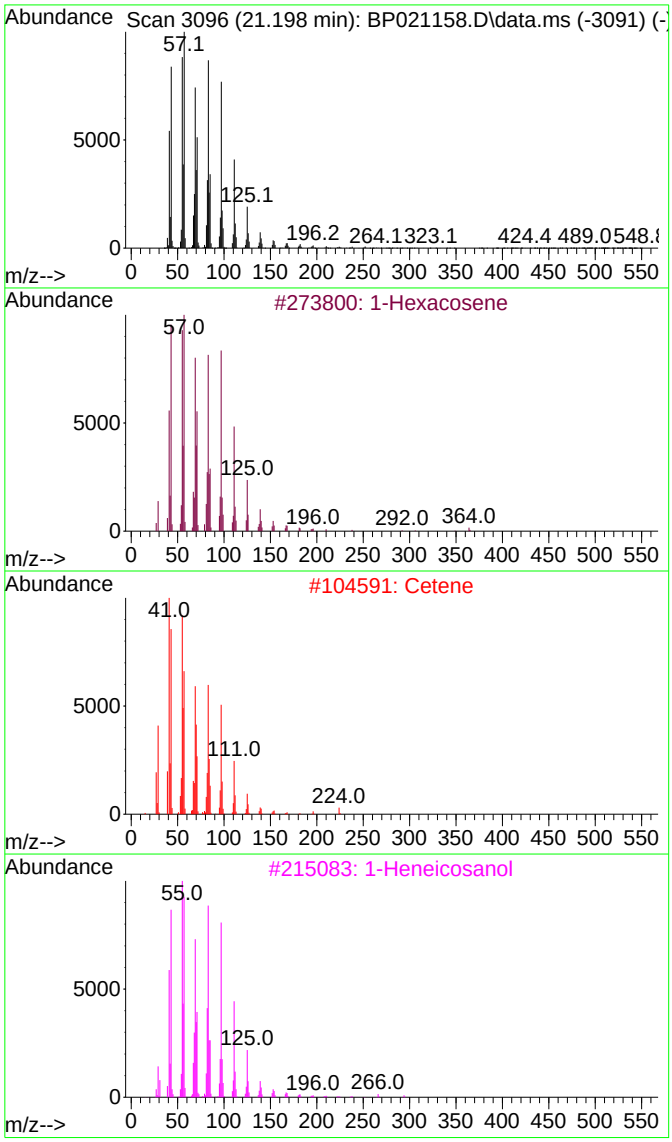
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

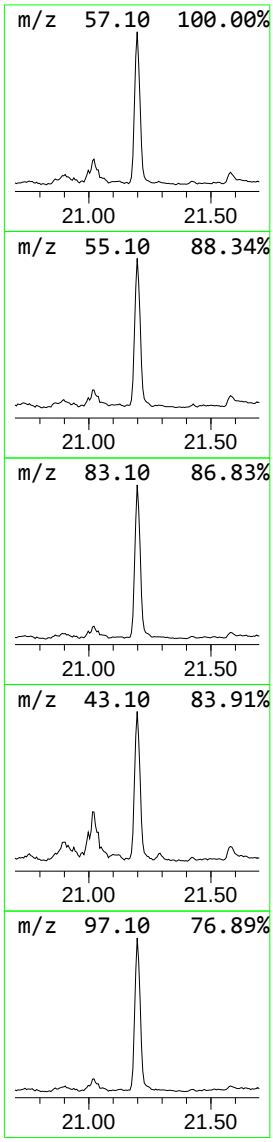
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.198	40.44 ng/ul	6589260	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	95
2	Cetene	224	C16H32	000629-73-2	95
3	1-Heneicosanol	312	C21H44O	015594-90-8	95
4	1-Tetracosene	336	C24H48	010192-32-2	94
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	94





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

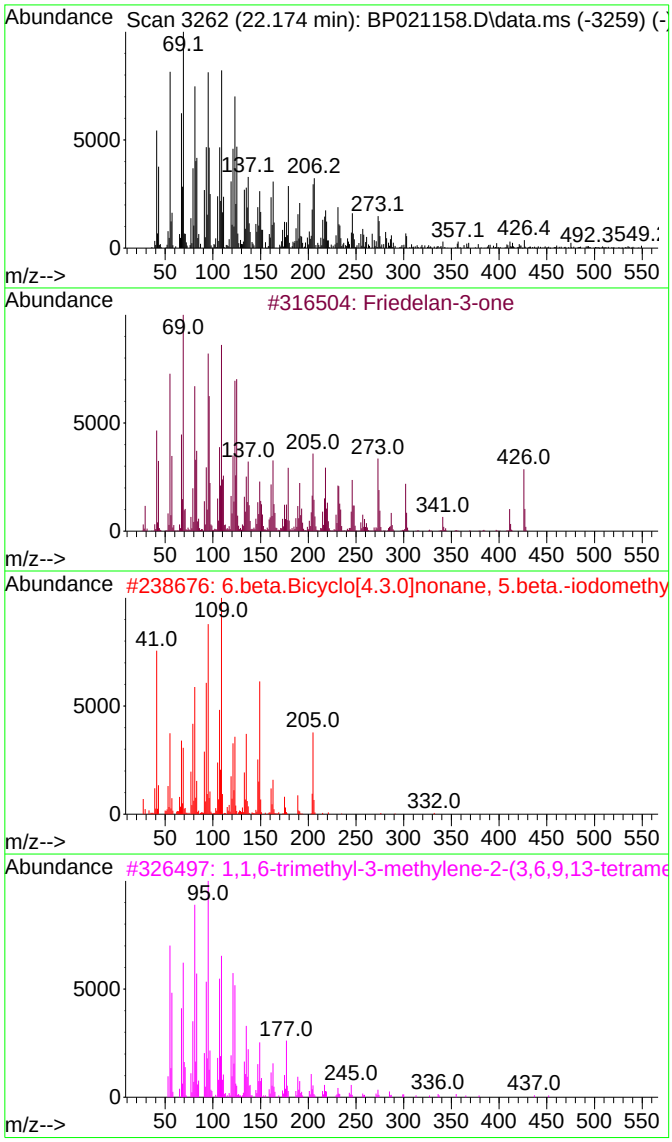
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

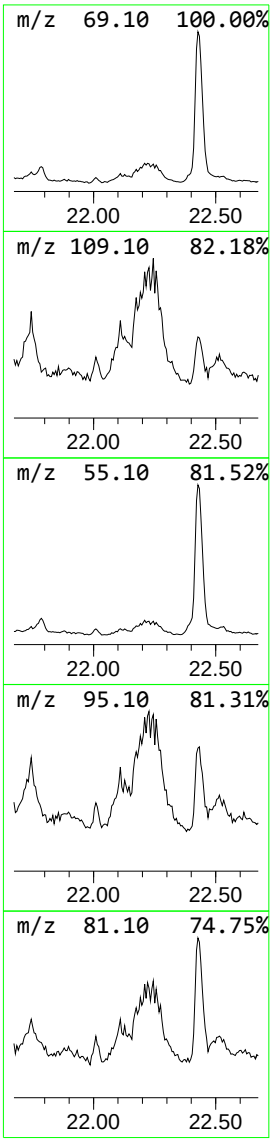
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	2.55 ng/ul	416035	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Friedelan-3-one	426	C30H50O	000559-74-0	89
2	6.beta.Bicyclo[4.3.0]nonane, 5.b...	332	C15H25I	1000195-85-9	84
3	1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	64
4	Friedelan-3-one	426	C30H50O	000559-74-0	62
5	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	59





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

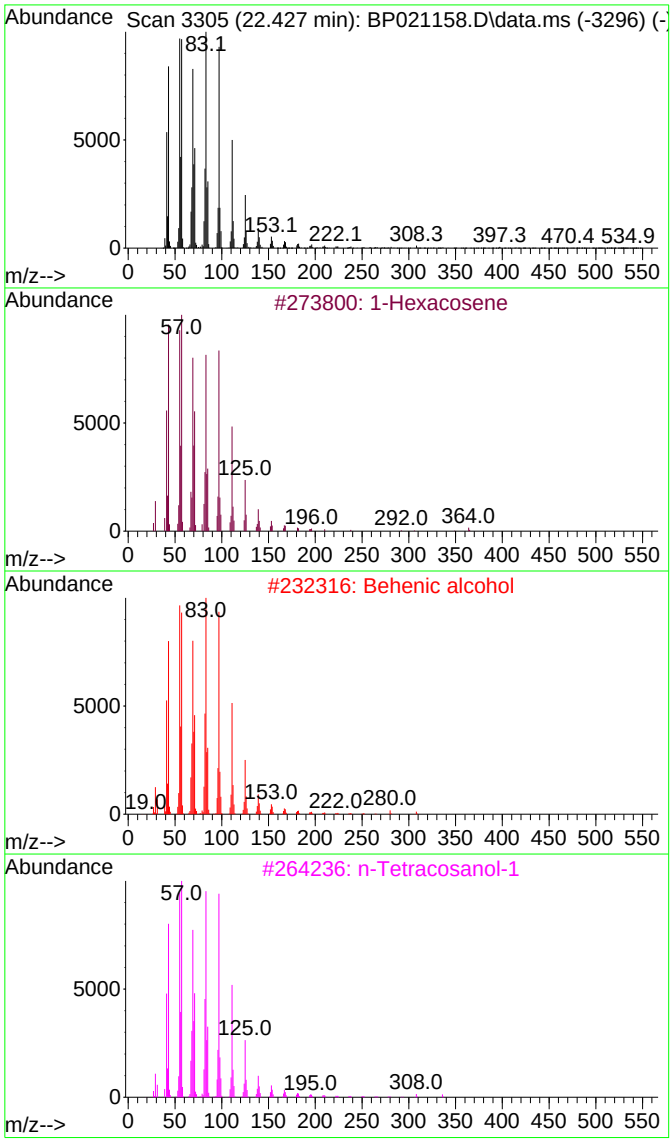
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

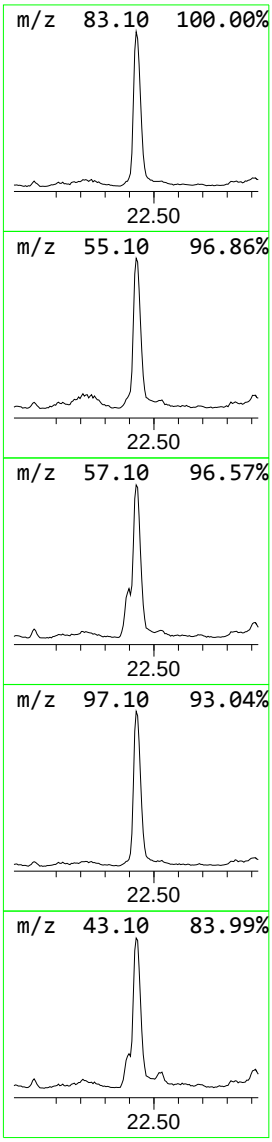
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.427	58.03 ng/ul	9455350	Chrysene-d12	21.551

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	95
2	Behenic alcohol	326	C22H46O	000661-19-8	95
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5	1-Heneicosanol	312	C21H44O	015594-90-8	94





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

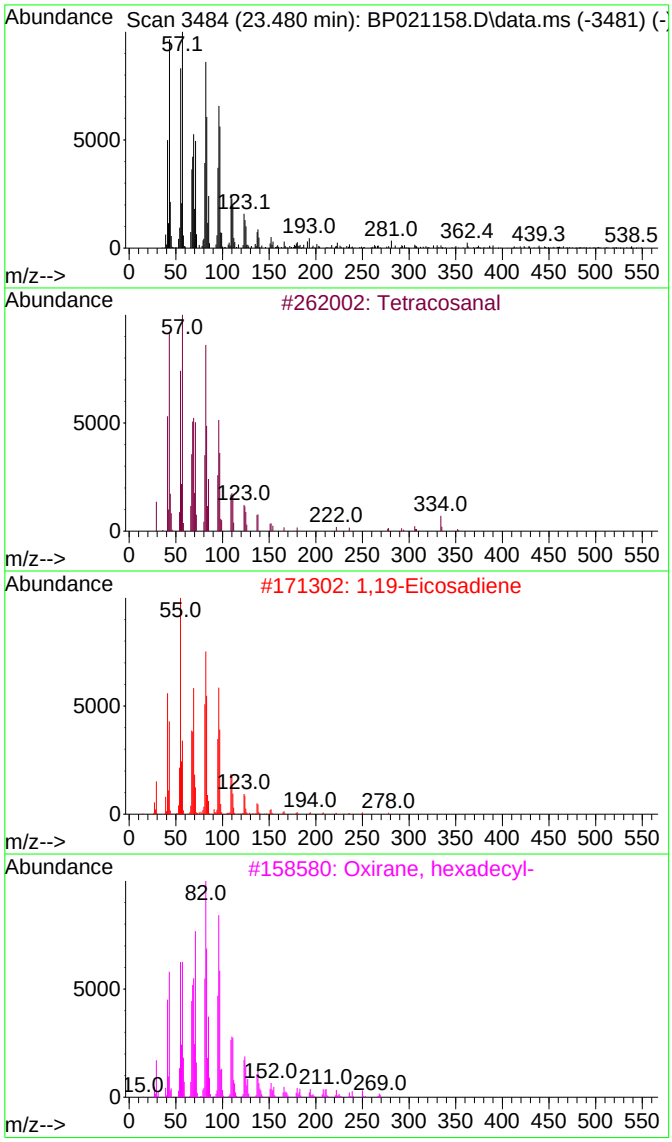
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

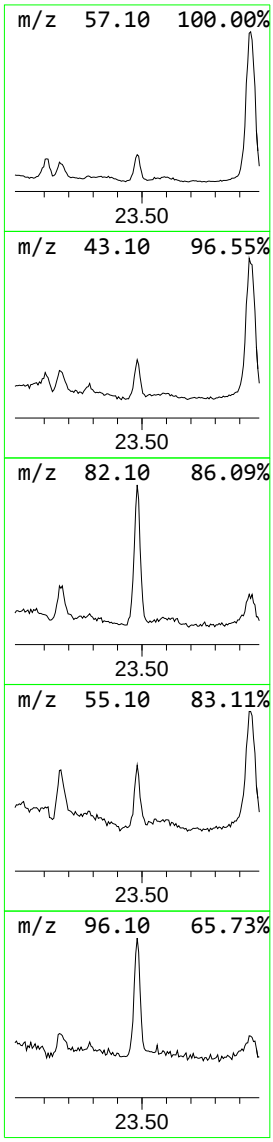
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Tetracosanal Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.480	4.51 ng/ul	692705	Perylene-d12	24.880

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosanal	352	C24H48O	057866-08-7	94
2	1,19-Eicosadiene	278	C20H38	014811-95-1	94
3	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
4	Hexacosanal	380	C26H52O	026627-85-0	94
5	Dotriacontanal	464	C32H64O	057878-00-9	94





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

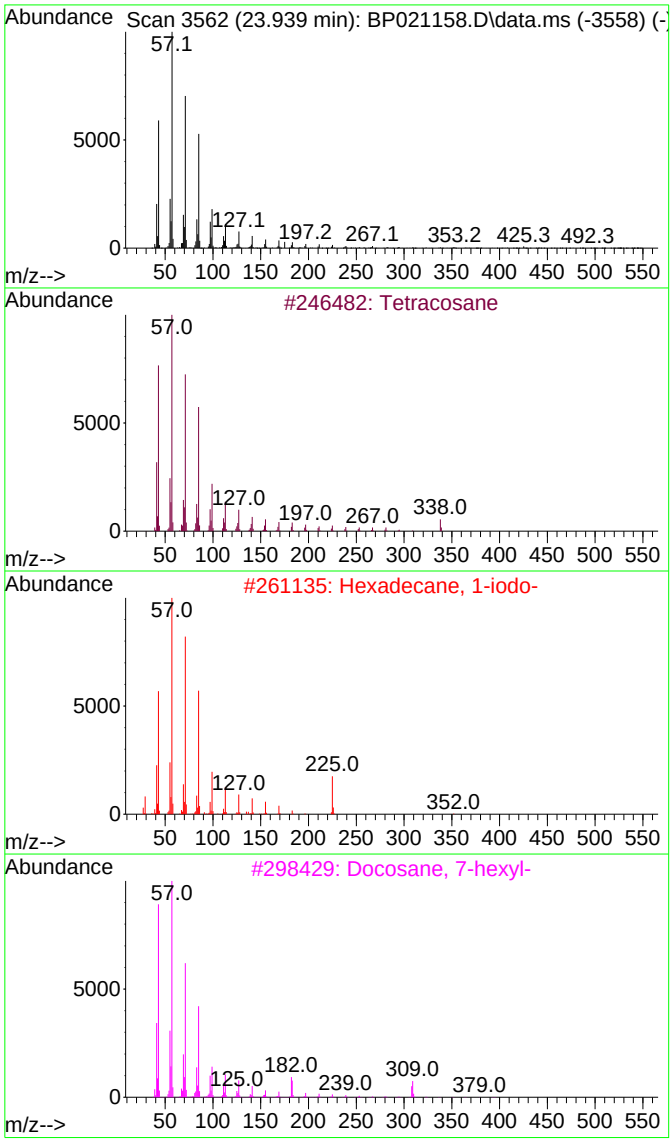
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

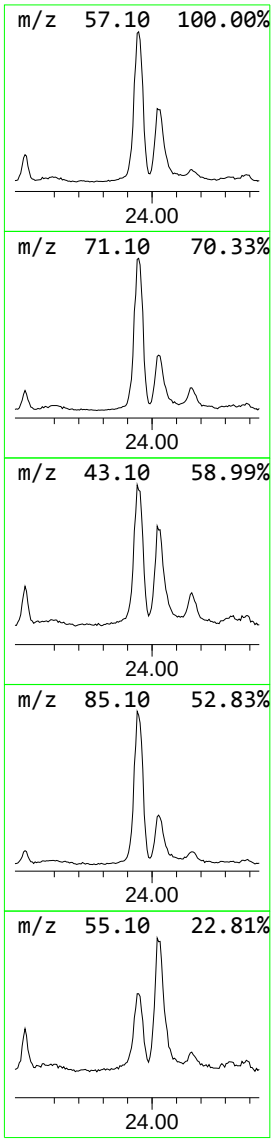
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.939	20.63 ng/ul	3170720	Perylene-d12	24.880

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
3	Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
4	Hexacosane	366	C26H54	000630-01-3	93
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	93





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Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

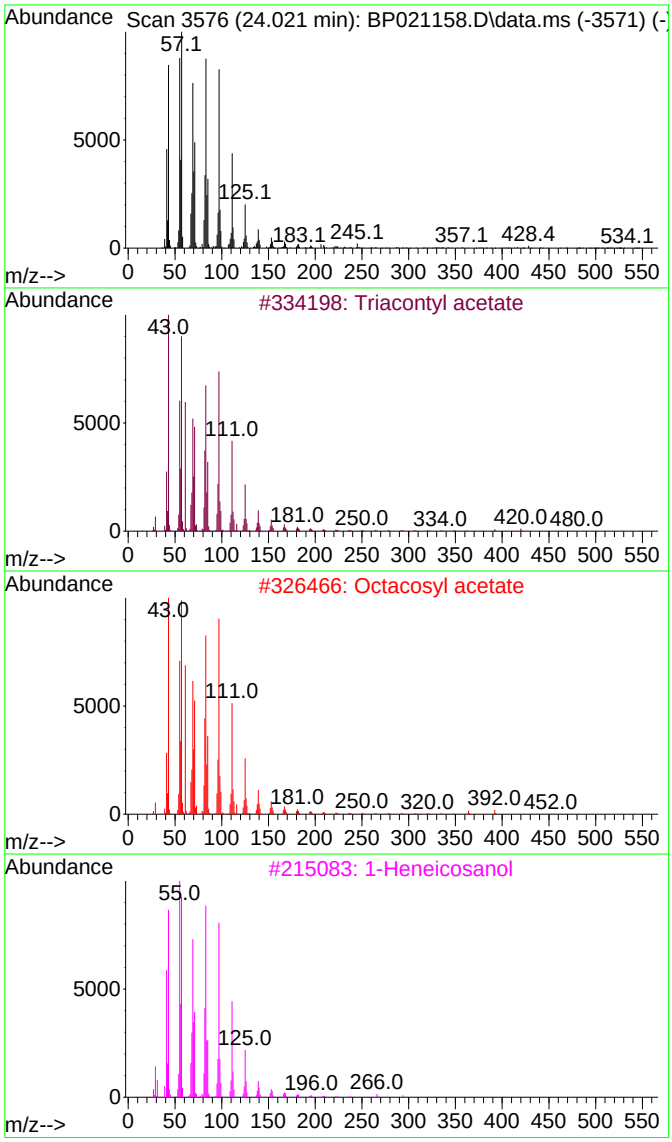
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

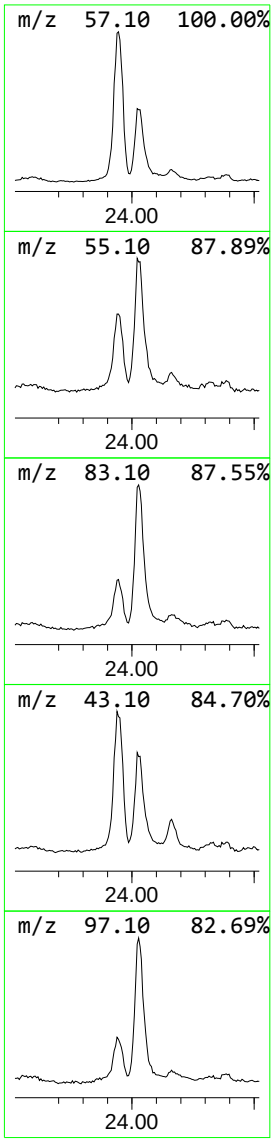
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Triacontyl acetate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.021	23.91 ng/ul	3675780	Perylene-d12	24.880

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontyl acetate	480	C32H64O2	041755-58-2	97
2	Octacosyl acetate	452	C30H60O2	018206-97-8	96
3	1-Heneicosanol	312	C21H44O	015594-90-8	95
4	1-Hexacosene	364	C26H52	018835-33-1	95
5	1-Tetracosene	336	C24H48	010192-32-2	94





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Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

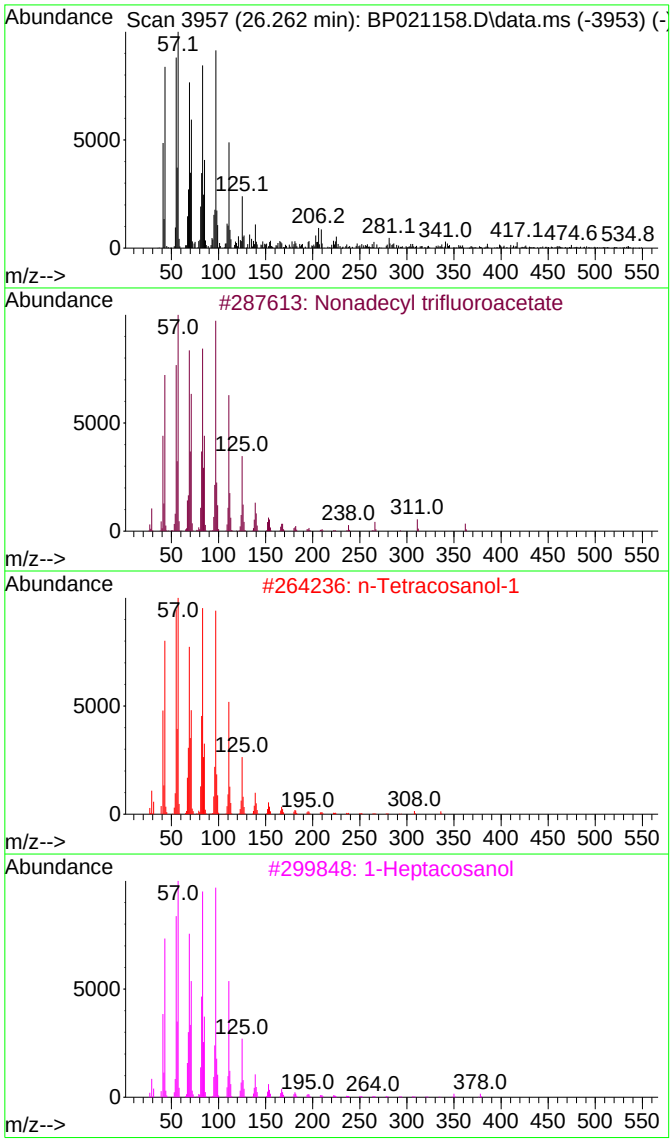
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

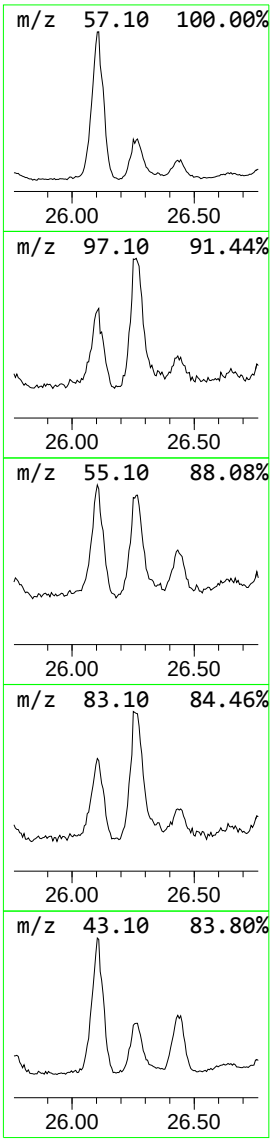
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 Nonadecyl trifluoroacetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.262	10.70 ng/ul	1644300	Perylene-d12	24.880

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	94
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
3	1-Heptacosanol	396	C27H56O	002004-39-9	91
4	1-Hexacosanol	382	C26H54O	000506-52-5	90
5	Octacosanol	410	C28H58O	000557-61-9	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021158.D
Acq On : 25 Jul 2024 22:26
Operator : MA/JU
Sample : P3263-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CG7

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
Isopropyl Alcohol	3.469	2.9	ng/ul	176947	1	7.651	1209460	20.0
Benzeneacetic acid	11.057	9.9	ng/ul	834762	2	10.422	1693450	20.0
1-Phenoxypropan...	11.181	3.8	ng/ul	323050	2	10.422	1693450	20.0
unknown-01	14.504	14.5	ng/ul	1474430	3	14.292	2032170	20.0
unknown-02	14.522	15.9	ng/ul	1610060	3	14.292	2032170	20.0
1-Dodecanol, 3,...	16.257	4.0	ng/ul	494595	4	17.110	2464160	20.0
(E)-4-(3-Hydrox...	16.504	6.4	ng/ul	786396	4	17.110	2464160	20.0
Stigmast-4-en-3...	17.751	3.3	ng/ul	404815	4	17.110	2464160	20.0
1,3-Hexadiene, ...	17.774	8.6	ng/ul	1063740	4	17.110	2464160	20.0
n-Hexadecanoic ...	18.039	18.8	ng/ul	2313070	4	17.110	2464160	20.0
(DEL) Alkane: S...	18.933	2.2	ng/ul	273557	4	17.110	2464160	20.0
2H-Azepin-2-one...	19.086	3.7	ng/ul	451924	4	17.110	2464160	20.0
Isopropyl linol...	19.221	2.5	ng/ul	304801	4	17.110	2464160	20.0
Amyl cyclopente...	19.292	6.4	ng/ul	783580	4	17.110	2464160	20.0
Octadecanoic acid	19.374	3.7	ng/ul	607562	5	21.551	3258650	20.0
(DEL) Alkane: S...	19.921	2.9	ng/ul	475868	5	21.551	3258650	20.0
(DEL) Alkane: S...	20.133	9.3	ng/ul	1509130	5	21.551	3258650	20.0
(DEL) Alkane: S...	20.680	2.3	ng/ul	370882	5	21.551	3258650	20.0
2-Tritriacontanone	20.998	2.4	ng/ul	398288	5	21.551	3258650	20.0
2-Nonacosanone	21.021	6.2	ng/ul	1005900	5	21.551	3258650	20.0
(DEL) Alkane: S...	21.198	40.4	ng/ul	6589260	5	21.551	3258650	20.0
5-Bromo-4-oxo-4...	22.174	2.5	ng/ul	416035	5	21.551	3258650	20.0
(DEL) Alkane: S...	22.427	58.0	ng/ul	9455350	5	21.551	3258650	20.0
Tetracosanal	23.480	4.5	ng/ul	692705	6	24.880	3074030	20.0
(DEL) Alkane: S...	23.939	20.6	ng/ul	3170720	6	24.880	3074030	20.0
Triacontyl acetate	24.021	23.9	ng/ul	3675780	6	24.880	3074030	20.0
Nonadecyl trifl...	26.262	10.7	ng/ul	1644300	6	24.880	3074030	20.0