LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCYD8

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-BP082324.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP021734.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	24	rVB	4720692	5707979	58.51%	3.331%
2	3.275	45	49	55	rVB	95115	115141	1.18%	0.067%
3	3.540	90	94	105	rBV	239962	336505	3.45%	0.196%
4	3.693	115	120	131	rBV	561579	831571	8.52%	0.485%
5	5.893	485	494	502	rVB2	59422	93048	0.95%	0.054%
6	6.499	590	597	605	rVB	288224	455138	4.67%	0.266%
7	6.799	643	648	651	rBV	59412	95217	0.98%	0.056%
8	6.840	651	655	664	rVV	96605	186156	1.91%	0.109%
9	6.963	666	676	692	rVV	1766783	2979047	30.54%	1.739%
10	7.116	696	702	719	rVB	1398081	2248715	23.05%	1.312%
11	7.328	730	738	745	rVV	1688885	2732977	28.01%	1.595%
12	7.393	745	749	759	rVV	100173	185339	1.90%	0.108%
13	7.616	782	787	796	rBV4	41398	84648	0.87%	0.049%
14	7.787	809	816	823	rBV	1660029	2762904	28.32%	1.612%
15	8.493	926	936	946	rBV	2197993	3644989	37.36%	2.127%
16	8.934	1004	1011	1023	rBV	1511148	2681060	27.48%	1.565%
17	9.493	1099	1106	1113	rBV	167375	266020	2.73%	0.155%
18	9.657	1127	1134	1141	rBV	1326400	2143145	21.97%	1.251%
19	9.846	1155	1166	1171	rBV	400623	981342	10.06%	0.573%
20	9.899	1172	1175	1188	rVB5	98808	210333	2.16%	0.123%
21	10.199	1217	1226	1248	rBV	1803339	3606242	36.96%	2.105%
22	10.575	1282	1290	1298	rBV	2341812	4023778	41.24%	2.348%
23	10.704	1304	1312	1333	rVB	985546	1852915	18.99%	1.081%
24	11.110	1376	1381	1389	rBV3	54580	109075	1.12%	0.064%
25	11.316	1410	1416	1435	rBV	296104	701890	7.19%	0.410%
26	12.169	1552	1561	1564	rBV2	36853	71064	0.73%	0.041%
27	12.216	1564	1569	1574	rVV3	67235	139862	1.43%	0.082%
28	12.640	1630	1641	1646	rBV7	30515	77348	0.79%	0.045%
29	12.793	1662	1667	1673	rVB2	93679	153635	1.57%	0.090%
30	13.181	1728	1733	1735	rBV4	71619	129086	1.32%	0.075%
31	13.310	1750	1755	1758	rBV2	183860	290900	2.98%	0.170%
32	13.528	1786	1792	1801	rBV	112899	205868	2.11%	0.120%
33	13.840	1838	1845	1852	rBV	3663401	5247132	53.78%	3.062%
34	13.904	1852	1856	1861	rVB	117293	158027	1.62%	0.092%
35	14.128	1884	1894	1902	rBV	3792673	6143386	62.97%	3.585%
36	14.204	1903	1907	1912	rVB2	66021	92596	0.95%	0.054%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

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-BP082324.MA.M
Title : SVOA CALIBRATION

37	14.287	1913	1921	1925	rBV3	240694	421423	4.32%	0.246%
38	14.404	1934	1941	1943	rVV	1699428	2419755	24.80%	1.412%
39	14.439	1943	1947	1953	rVV	3250237	5270209	54.02%	3.076%
40	14.522	1956	1961	1974	rVB2	262337	535325	5.49%	0.312%
41	14.645	1974	1982	1988	rBV	1508658	3040137	31.16%	1.774%
42	14.698	1988	1991	1999	rVB2	437068	748703	7.67%	0.437%
43	14.775	1999	2004	2012	rVB2	189185	432738	4.44%	0.253%
44	14.875	2015	2021	2024	rBV3	132489	214343	2.20%	0.125%
45	14.969	2034	2037	2047	rVB	51261	98929	1.01%	0.058%
46	15.110	2057	2061	2067	rVB6	48722	82044	0.84%	0.048%
47	15.251	2081	2085	2091	rVB3	59447	100996	1.04%	0.059%
48	15.445	2111	2118	2128	rVB	4337070	7403924	75.89%	4.321%
49	15.557	2130	2137	2144	rBV	1330280	1977067	20.27%	1.154%
50	15.692	2156	2160	2167	rVB6	64240	114068	1.17%	0.067%
51	15.798	2173	2178	2182	rBV2	119375	210029	2.15%	0.123%
52	15.845	2183	2186	2190	rVV2	79918	108165	1.11%	0.063%
53	15.945	2193	2203	2208	rVB2	453801	950724	9.74%	0.555%
54	16.034	2210	2218	2224	rBV3	265289	444449	4.56%	0.259%
55	16.092	2224	2228	2237	rBV6	63210	152757	1.57%	0.089%
56	16.192	2241	2245	2250	rBV2	173210	244617	2.51%	0.143%
57	16.345	2268	2271	2275	rBV4	57096	82973	0.85%	0.048%
58	16.616	2311	2317	2329	rVB4	369285	717686	7.36%	0.419%
59	16.728	2332	2336	2344	rBV	78033	140654	1.44%	0.082%
60	16.963	2372	2376	2381	rVV5	43597	84192	0.86%	0.049%
61	17.051	2382	2391	2398	rVB2	168819	358210	3.67%	0.209%
62	17.192	2410	2415	2417	rBV2	108917	168270	1.72%	0.098%
63	17.239	2417	2423	2430	rVV	3993647	6648687	68.15%	3.880%
64	17.339	2434	2440	2451	rVB2	5793577	8121284	83.24%	4.740%
65	17.451	2453	2459	2463	rBV4	84743	165404	1.70%	0.097%
66	17.551	2470	2476	2480	rVB5	109193	224344	2.30%	0.131%
67	17.645	2487	2492	2502	rVB	302724	554828	5.69%	0.324%
68	17.910	2532	2537	2540	rBV	179143	289072	2.96%	0.169%
69	17.951	2541	2544	2553	rVB4	158098	280726	2.88%	0.164%
70	18.116	2569	2572	2575	rBV4	79593	107777	1.10%	0.063%
71	18.180	2578	2583	2596	rBV	1871197	3329007	34.12%	1.943%
72	18.310	2601	2605	2609	rVB2	219270	292716	3.00%	0.171%
73	18.486	2632	2635	2638	rBV3	76465	106971	1.10%	0.062%
74	18.533	2640	2643	2648	rVB2	138226	213287	2.19%	0.124%
75	18.692	2658	2670	2677	rBV2	202429	762387	7.81%	0.445%
76	18.827	2690	2693	2696	rBV3	86588	122453	1.26%	0.071%

LSC Area Percent Report

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Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

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-BP082324.MA.M
Title : SVOA CALIBRATION

77	19.075	2730	2735	2740	rBV3	560789	1052741	10.79%	0.614%
78	19.116	2740	2742	2747	rVB3	267511	368877	3.78%	0.215%
79	19.245	2761	2764	2765	rBV	112064	133338	1.37%	0.078%
80	19.357	2779	2783	2785	rBV2	263125	378317	3.88%	0.221%
81	19.386	2785	2788	2803	rVB3	662678	1502202	15.40%	0.877%
82	19.510	2804	2809	2819	rVB	873132	1445325	14.81%	0.843%
83	19.710	2837	2843	2852	rBV	5793096	9756049	100.00%	5.694%
84	20.086	2904	2907	2911	rBV2	740402	987508	10.12%	0.576%
85	20.127	2911	2914	2919	rVB	831051	1125091	11.53%	0.657%
86	20.274	2935	2939	2943	rBV2	608399	946291	9.70%	0.552%
87	20.598	2991	2994	2997	rBV2	450546	608364	6.24%	0.355%
88	20.633	2997	3000	3009	rVB	1548766	2025436	20.76%	1.182%
89	20.739	3016	3018	3023	rVB	469689	534588	5.48%	0.312%
90	21.016	3062	3065	3068	rBV	369544	397793	4.08%	0.232%
91	21.145	3083	3087	3094	rVB	7514725	9648135	98.89%	5.631%
92	21.221	3098	3100	3107	rVB3	497843	614769	6.30%	0.359%
93	21.310	3110	3115	3119	rVV	5039895	7545075	77.34%	4.403%
94	21.351	3119	3122	3131	rVB	2009730	3215243	32.96%	1.876%
95	21.686	3173	3179	3185	rBV2	3352957	6190295	63.45%	3.613%
96	21.751	3186	3190	3197	rVB	859289	1420758	14.56%	0.829%
97	22.616	3332	3337	3356	rVB	1647810	3428451	35.14%	2.001%
98	24.862	3710	3719	3727	rVB	3418839	7907040	81.05%	4.614%
99	25.086	3749	3757	3776	rVB	2275595	6982581	71.57%	4.075%
100	26.627	4013	4019	4035	rVB2	1116572	3652951	37.44%	2.132%

Sum of corrected areas: 171352626

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
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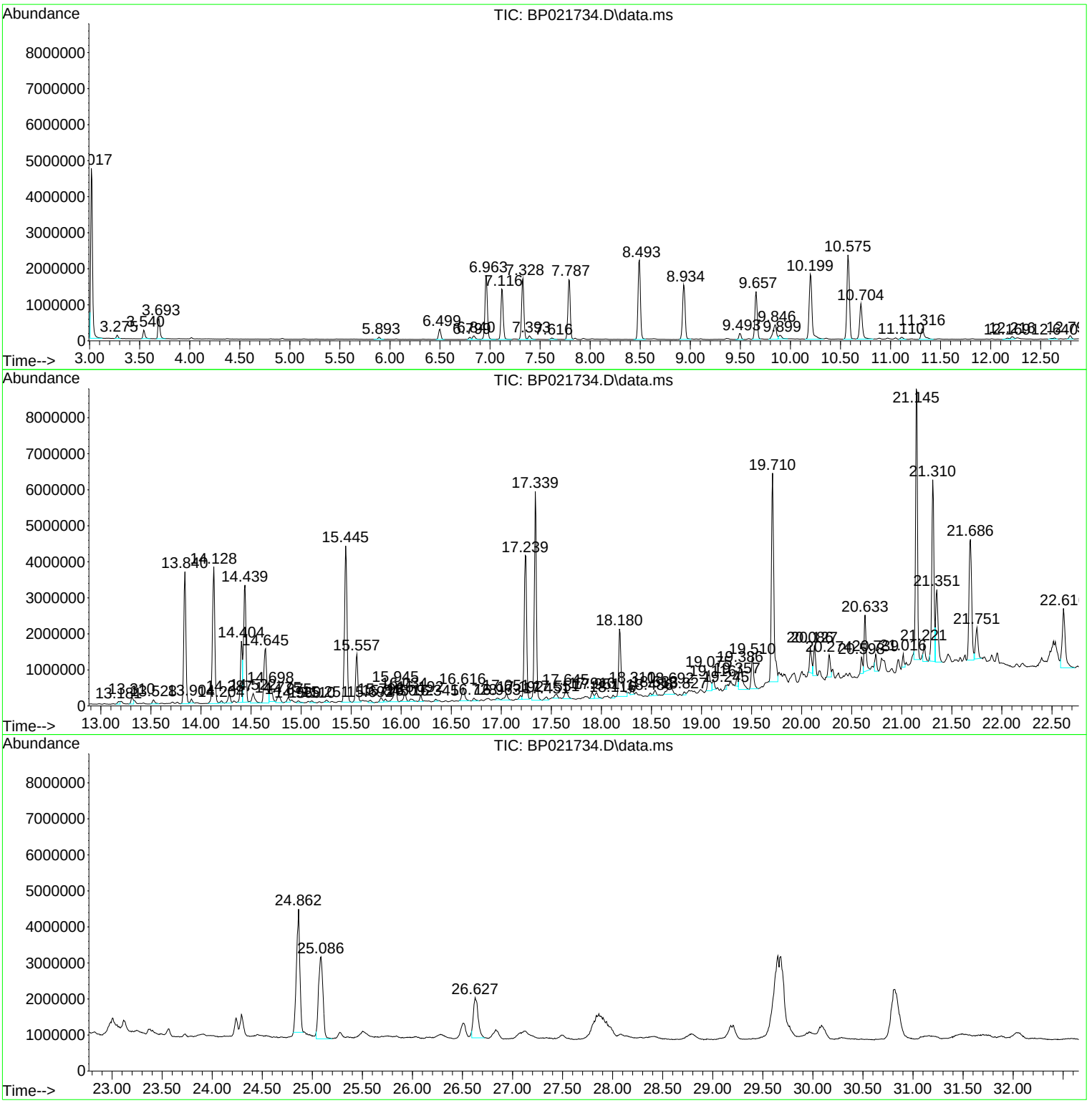
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.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\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

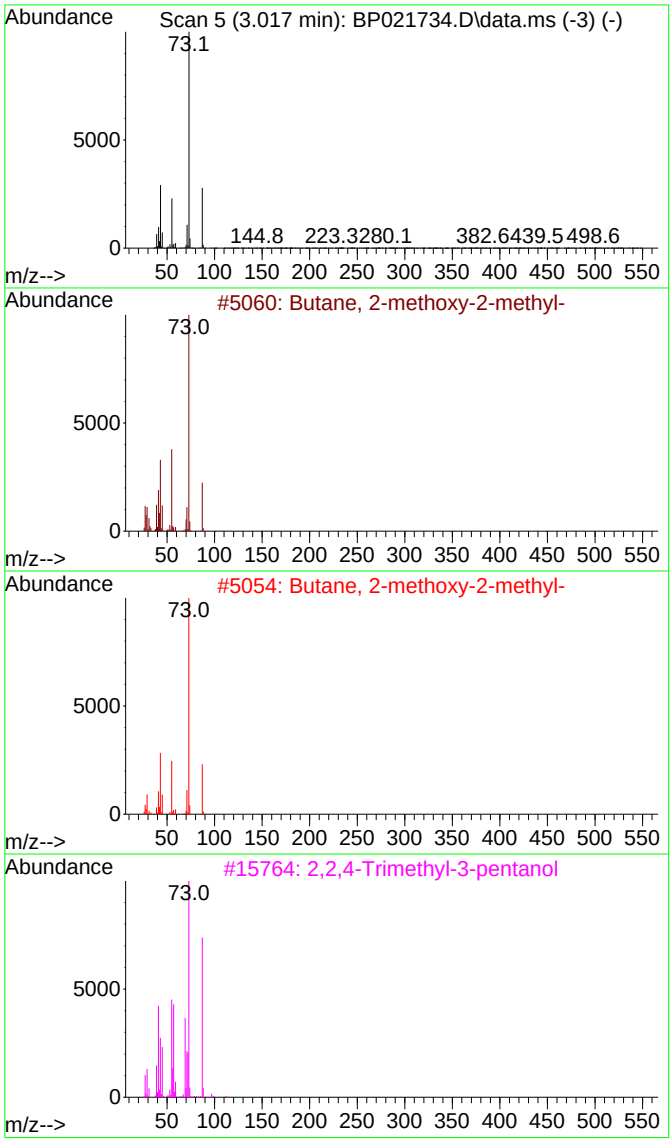
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

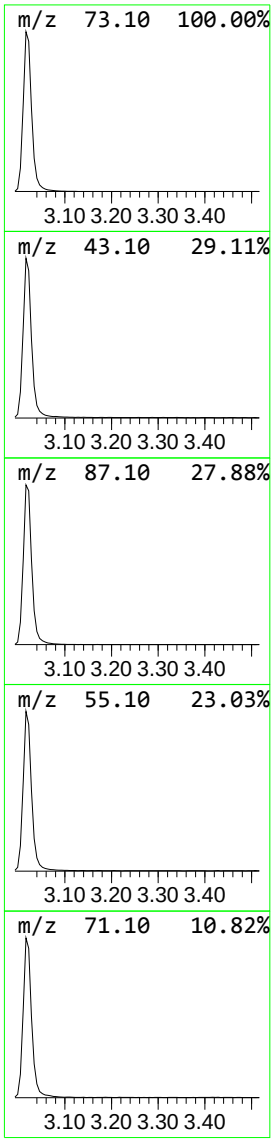
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	41.32 ng/ul	5707980	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

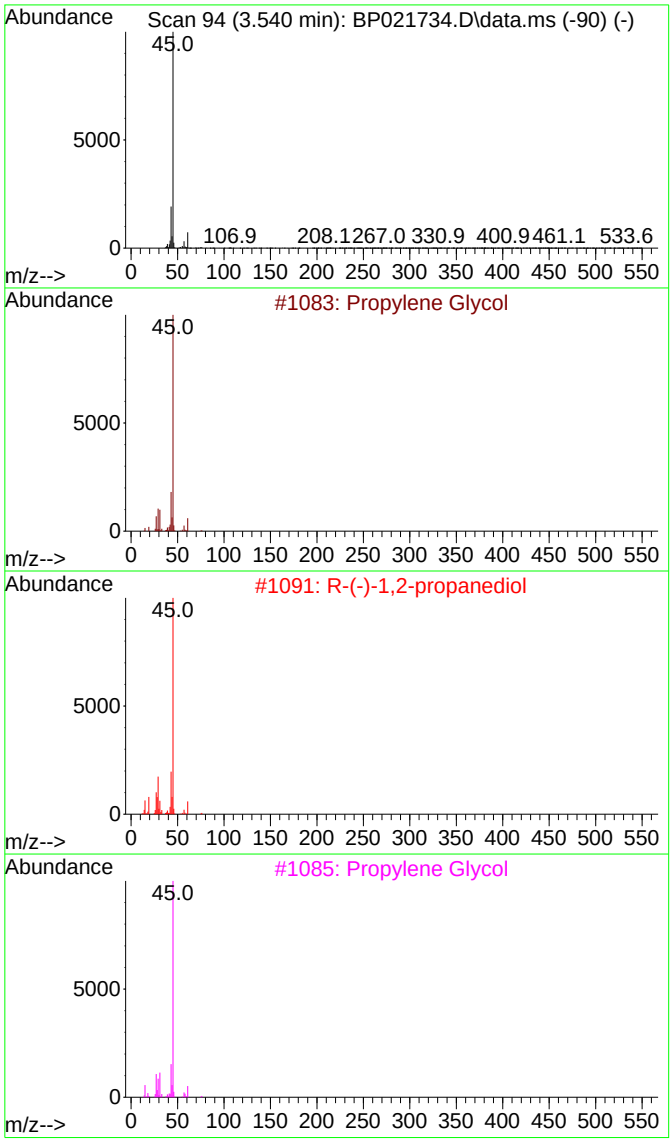
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

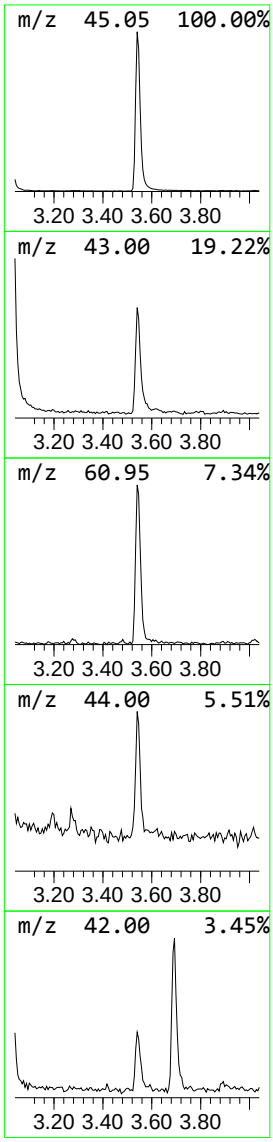
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Propylene Glycol Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.540	2.44 ng/ul	336505	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	90
2	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3	Propylene Glycol	76	C3H8O2	000057-55-6	78
4	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
5	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

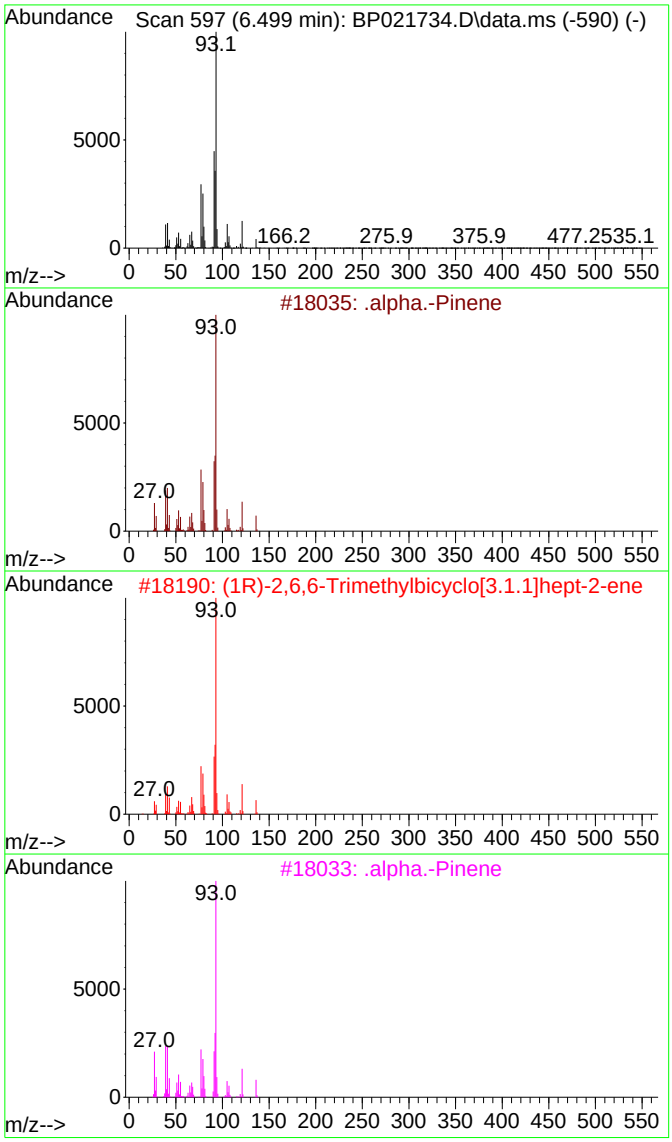
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

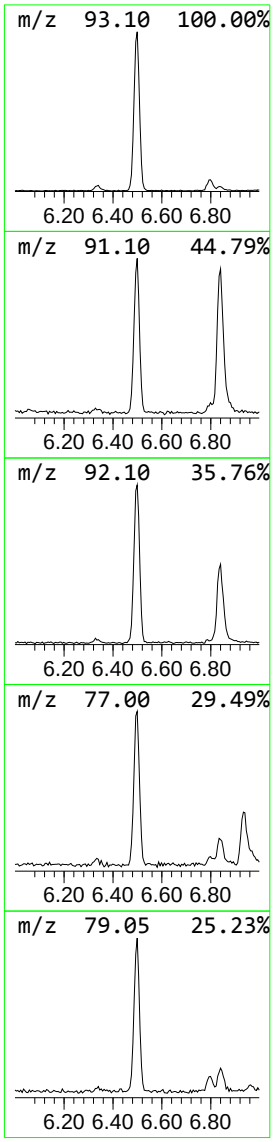
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 .alpha.-Pinene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.499	3.29 ng/ul	455138	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	94
2	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	94
3	.alpha.-Pinene	136	C10H16	000080-56-8	91
4	.alpha.-Pinene	136	C10H16	000080-56-8	91
5	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

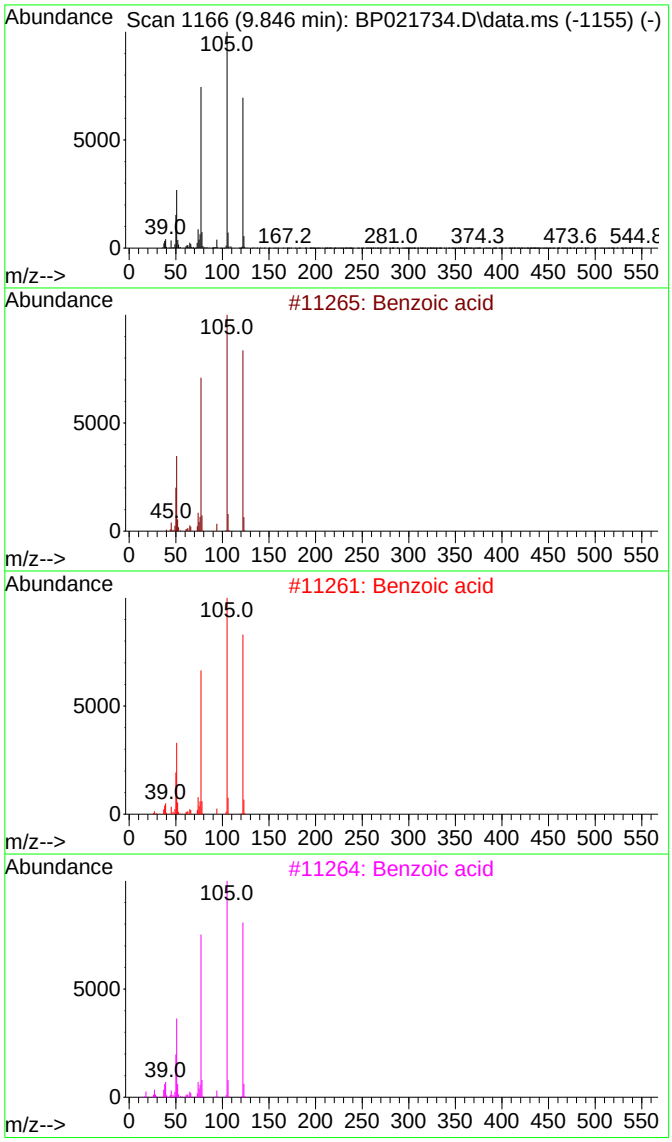
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

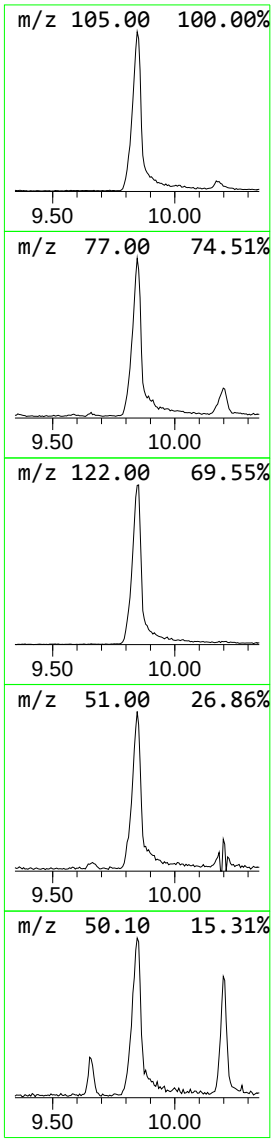
****

Peak Number 4 Benzoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.846	4.88 ng/ul	981342	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid	122	C7H6O2	000065-85-0	96
2	Benzoic acid	122	C7H6O2	000065-85-0	94
3	Benzoic acid	122	C7H6O2	000065-85-0	94
4	Benzoic acid	122	C7H6O2	000065-85-0	91
5	Benzoic acid, silver(1+) salt	228	C7H5AgO2	000532-31-0	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

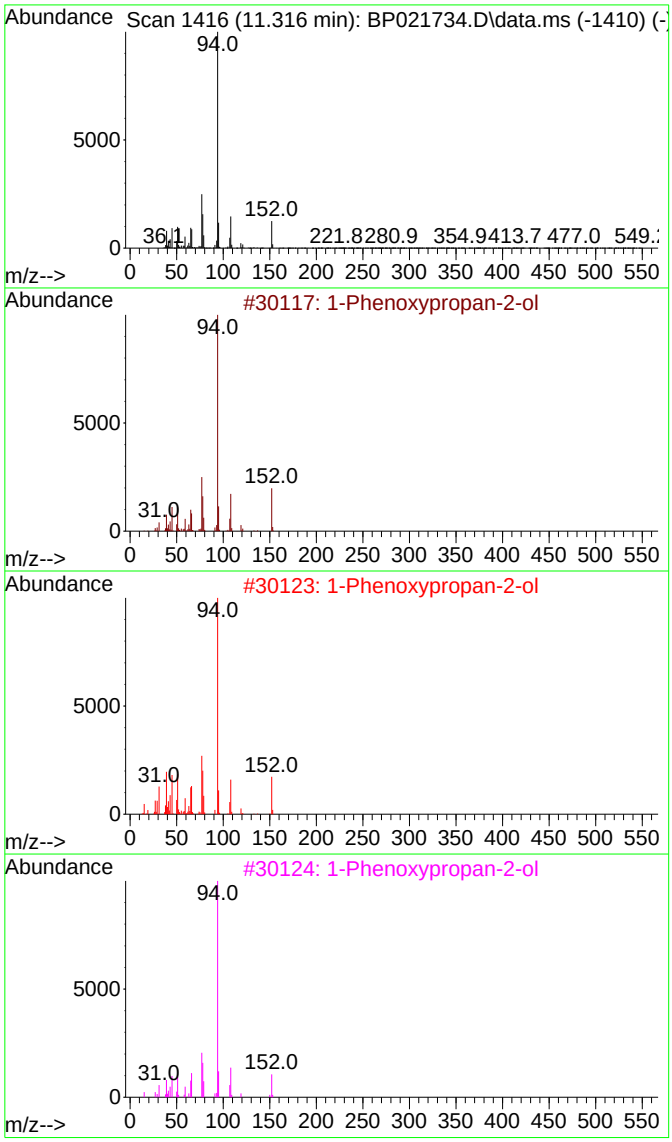
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

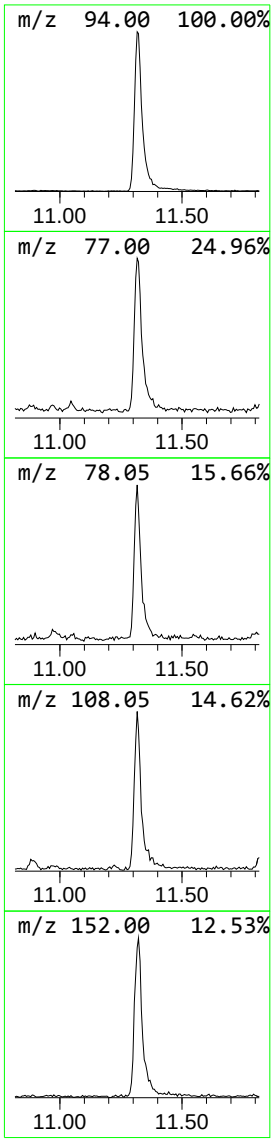
****

Peak Number 5 1-Phenoxypropan-2-ol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.316	3.49 ng/ul	701890	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	90
5	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87





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Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

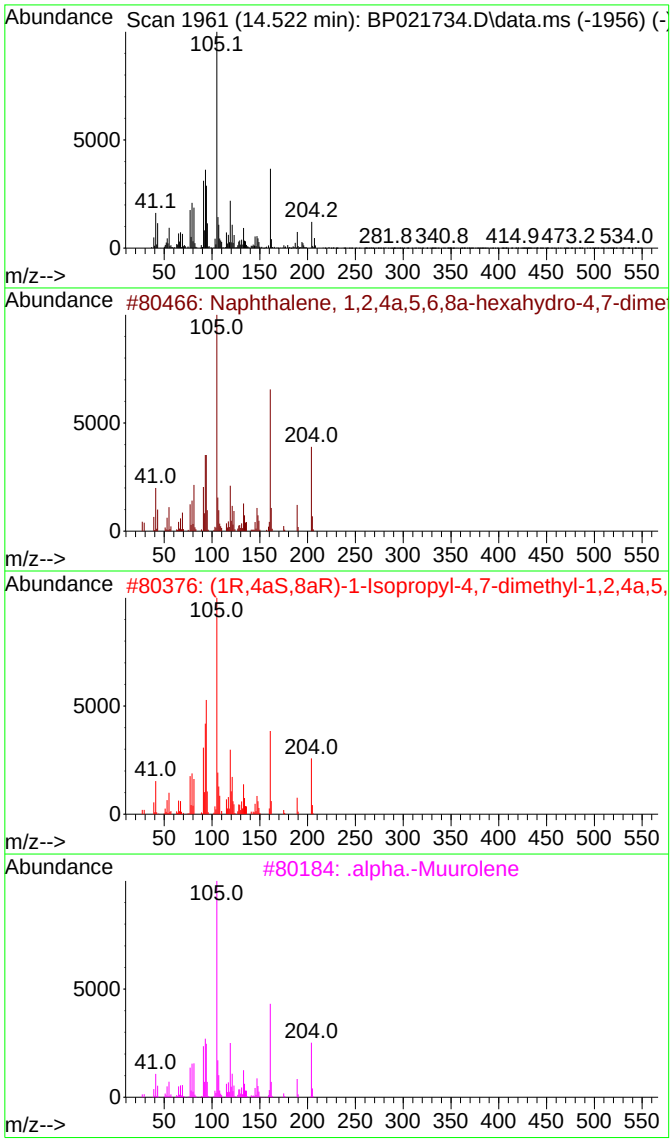
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

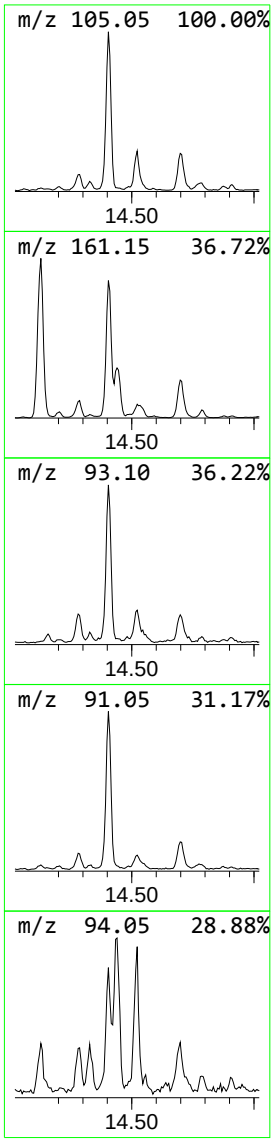
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Naphthalene, 1,2,4a,5,6,8a-... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.522	2.03 ng/ul	535325	Acenaphthene-d10	14.440

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,4a,5,6,8a-hexah...	204	C15H24	031983-22-9	93
2	(1R,4aS,8aR)-1-Isopropyl-4,7-dim...	204	C15H24	020085-19-2	93
3	.alpha.-Muurolene	204	C15H24	010208-80-7	93
4	Naphthalene, 1,2,4a,5,6,8a-hexah...	204	C15H24	000483-75-0	91
5	Naphthalene, 1,2,4a,5,6,8a-hexah...	204	C15H24	031983-22-9	91





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Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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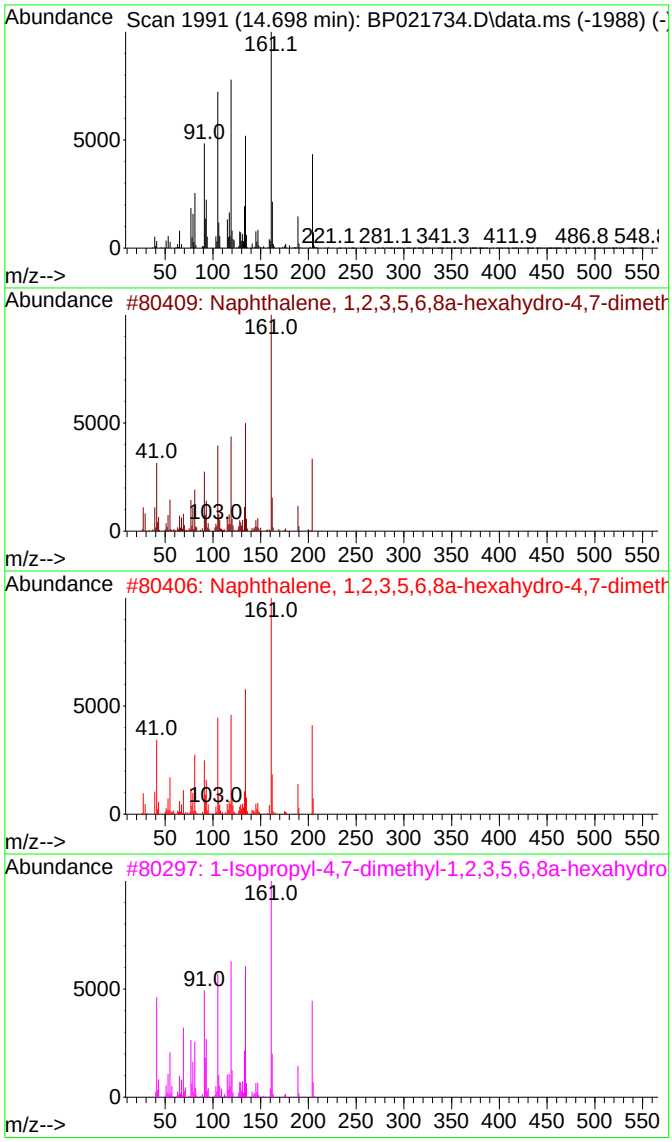
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

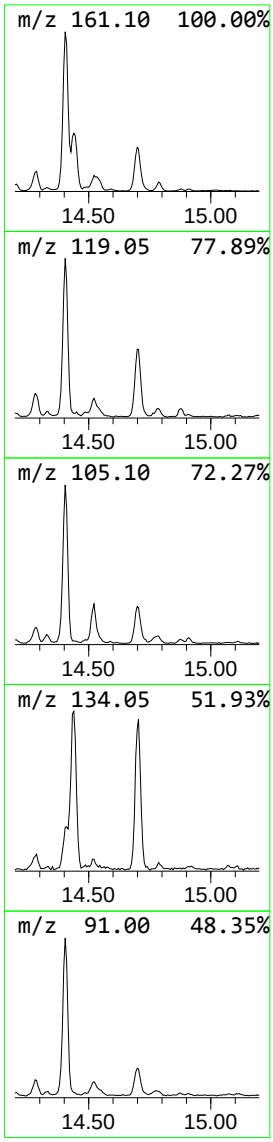
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Naphthalene, 1,2,3,5,6,8a-h... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.698	2.84 ng/ul	748703	Acenaphthene-d10	14.440

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,5,6,8a-hexahy...	204	C15H24	000483-76-1	97
2	Naphthalene, 1,2,3,5,6,8a-hexahy...	204	C15H24	000483-76-1	97
3	1-Isopropyl-4,7-dimethyl-1,2,3,5...	204	C15H24	016729-01-4	97
4	Naphthalene, 1,2,3,5,6,8a-hexahy...	204	C15H24	000483-76-1	93
5	Naphthalene, 1,2,4a,5,8,8a-hexah...	204	C15H24	000523-47-7	91





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Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

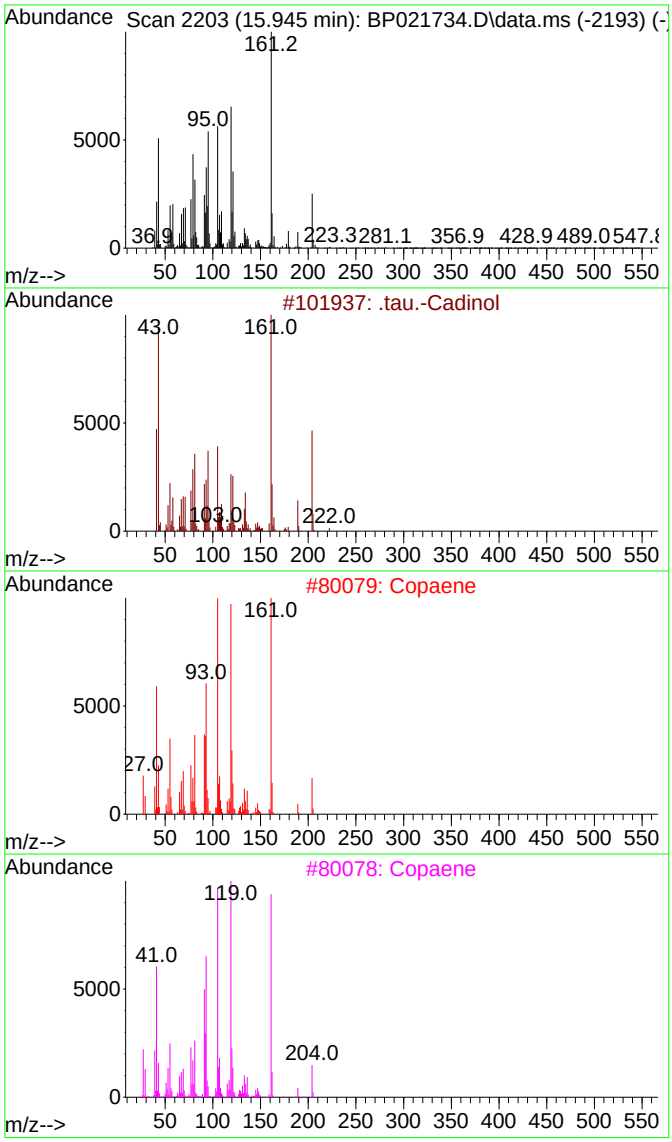
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

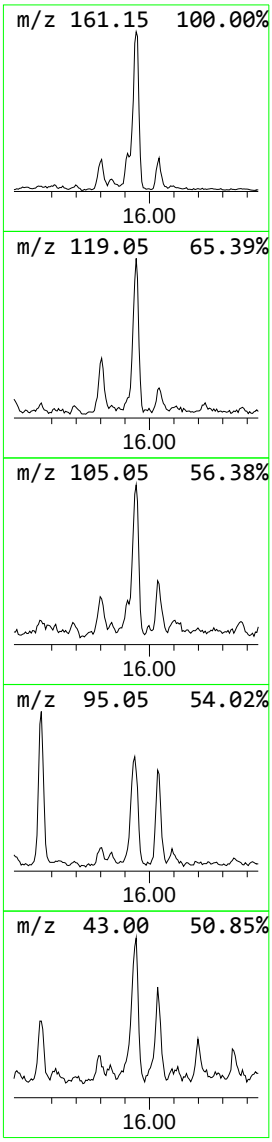
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 .tau.-Cadinol Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	2.86 ng/ul	950724	Phenanthrene-d10	17.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.tau.-Cadinol	222	C15H26O	005937-11-1	93
2	Copaene	204	C15H24	003856-25-5	91
3	Copaene	204	C15H24	003856-25-5	91
4	.alpha.-Cubebene	204	C15H24	017699-14-8	90
5	.gamma.-Muurolene	204	C15H24	030021-74-0	89





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

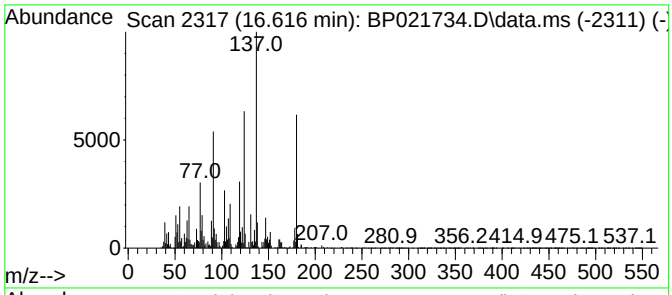
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

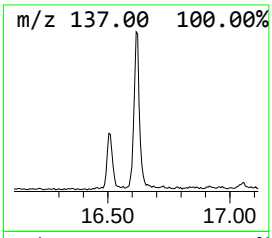
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

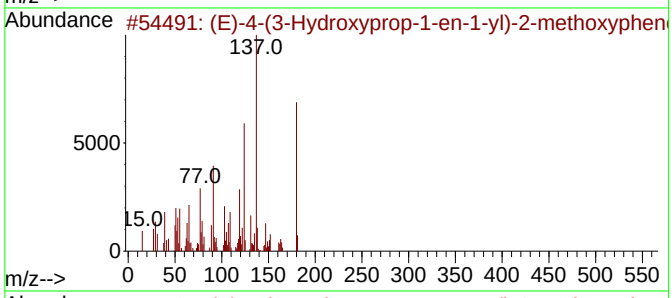
****
Peak Number 9 (E)-4-(3-Hydroxyprop-1-en-1... Concentration Rank 23

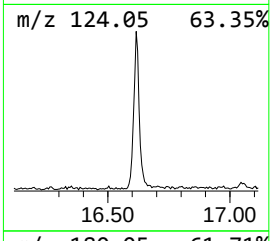
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.616	2.16 ng/ul	717686	Phenanthrene-d10	17.239

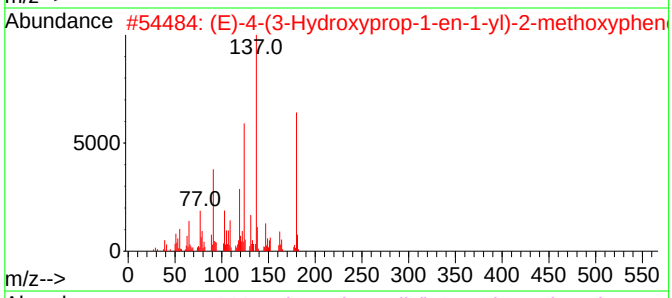
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-4-(3-Hydroxyprop-1-en-1-yl)-...	180	C10H12O3	032811-40-8	98
2	(E)-4-(3-Hydroxyprop-1-en-1-yl)-...	180	C10H12O3	032811-40-8	94
3	4-(1-Hydroxyallyl)-2-methoxyphenol	180	C10H12O3	112465-50-6	64
4	2,3,4,5-Tetrafluorobenzyl alcohol	180	C7H4F4O	053072-18-7	56
5	3,7-Benzofurandiol, 2,3-dihydro-...	180	C10H12O3	017781-15-6	49

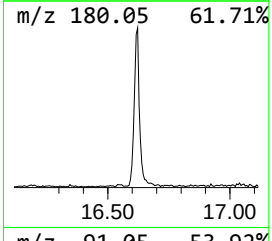


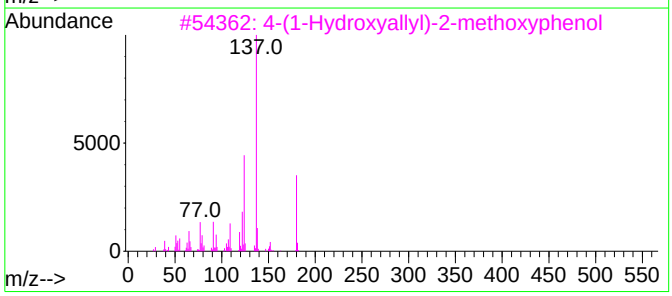


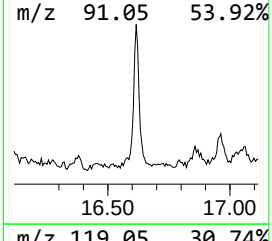


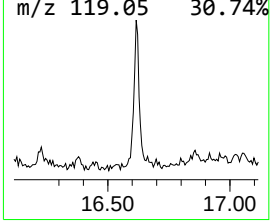












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

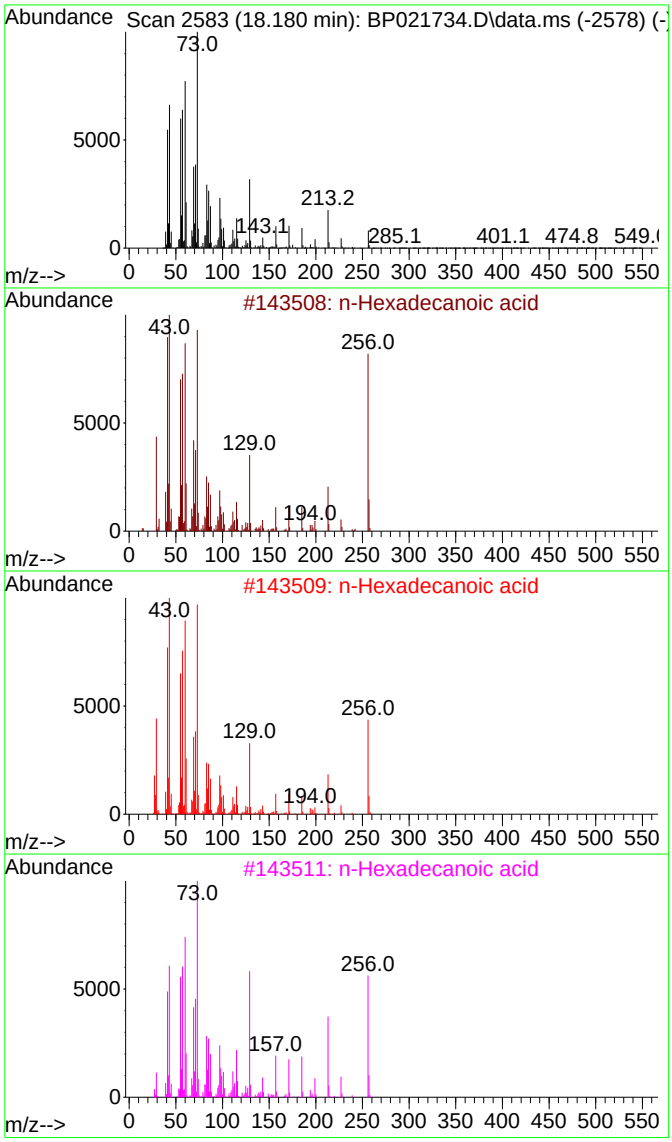
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

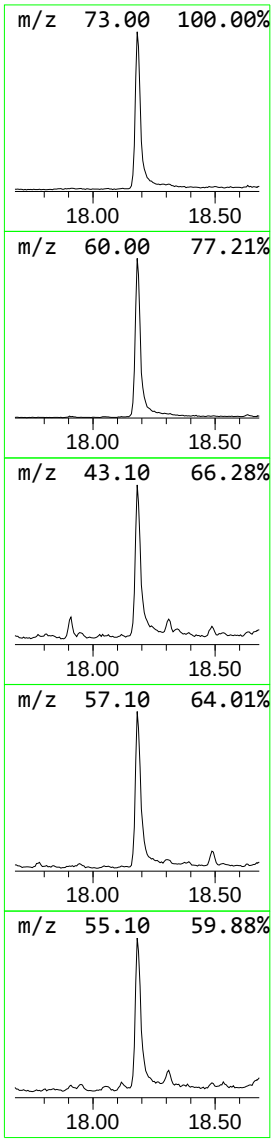
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	10.01 ng/ul	3329010	Phenanthrene-d10	17.239

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	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

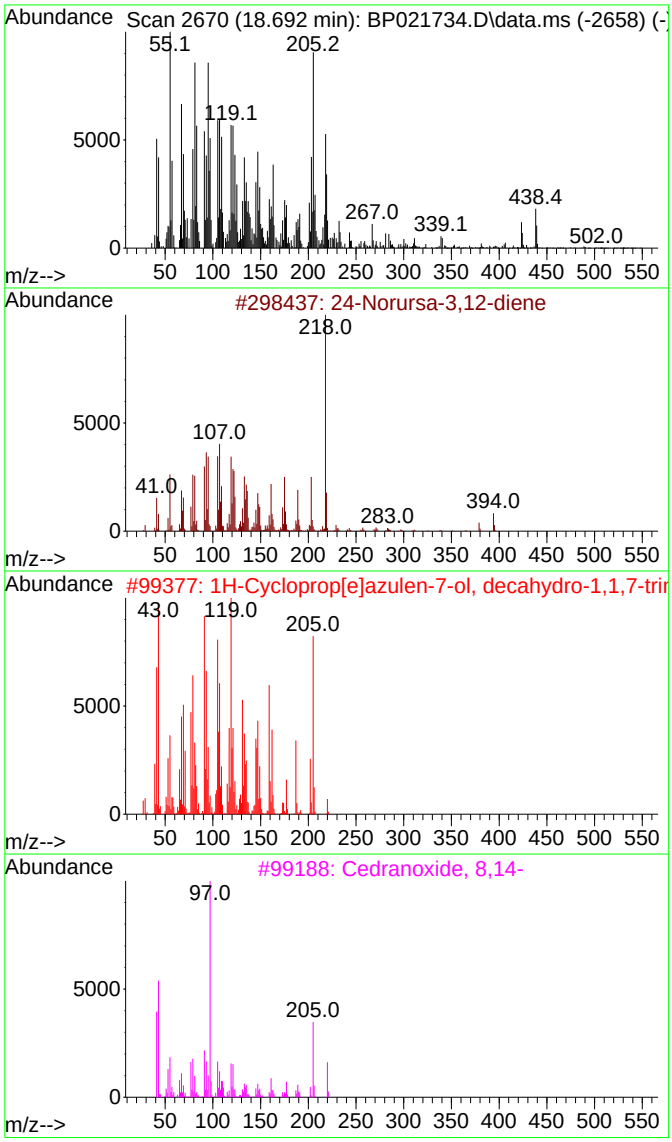
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

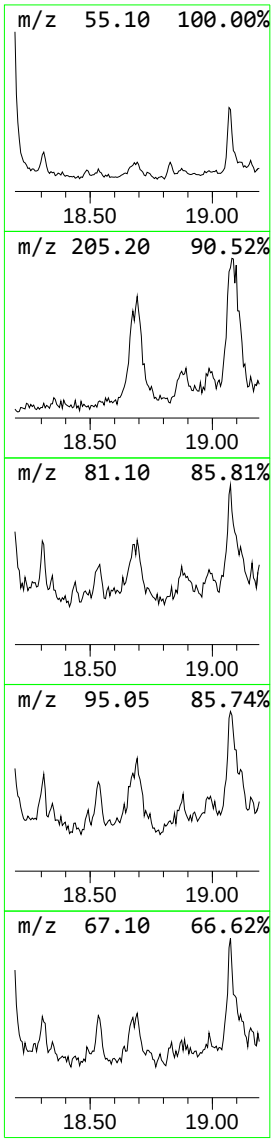
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 24-Norursa-3,12-diene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.692	2.29 ng/ul	762387	Phenanthrene-d10	17.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	24-Norursa-3,12-diene	394	C29H46	201358-25-0	70
2	1H-Cycloprop[e]azulen-7-ol, deca...	220	C15H24O	006750-60-3	55
3	Cedranoxide, 8,14-	220	C15H24O	018319-31-8	45
4	(R)-2-((4aS,8aR)-4a-Methyl-8-met...	220	C15H24O	028102-68-3	43
5	Lup-20(29)-en-3-one	424	C30H48O	001617-70-5	35





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

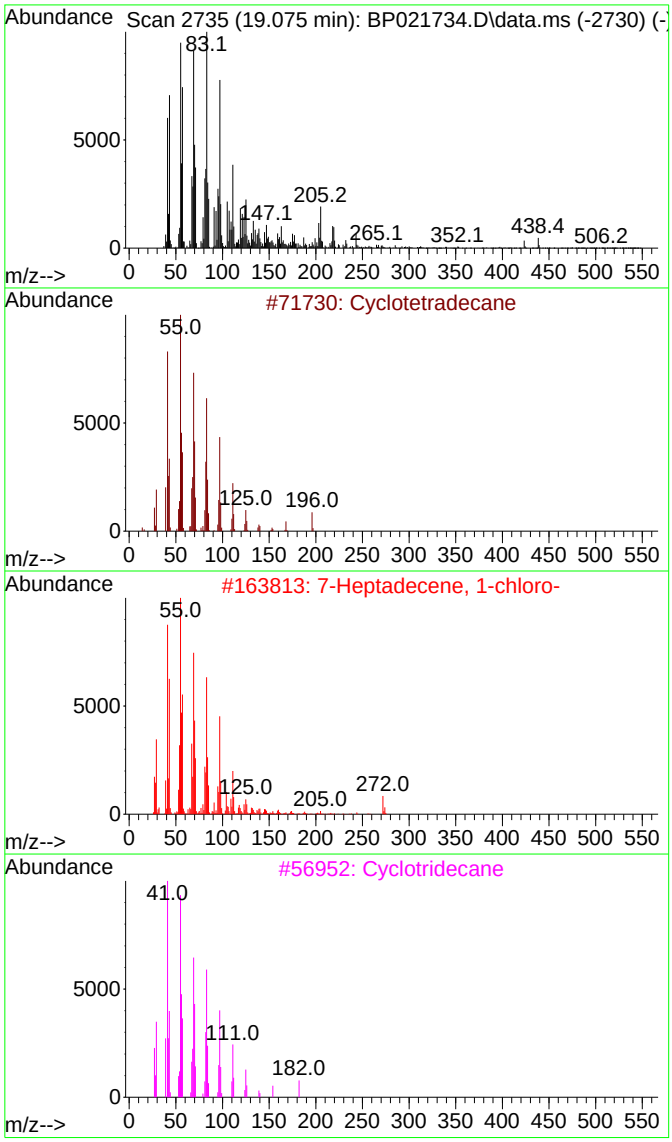
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

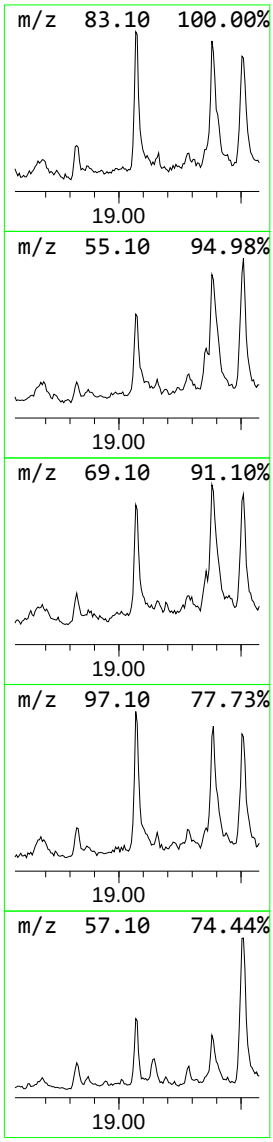
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Cyclic19.075 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.075	3.17 ng/ul	1052740	Phenanthrene-d10	17.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	94
2	7-Heptadecene, 1-chloro-	272	C17H33Cl	056554-78-0	93
3	Cyclotridecane	182	C13H26	000295-02-3	93
4	Cyclohexadecane	224	C16H32	000295-65-8	92
5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	90





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

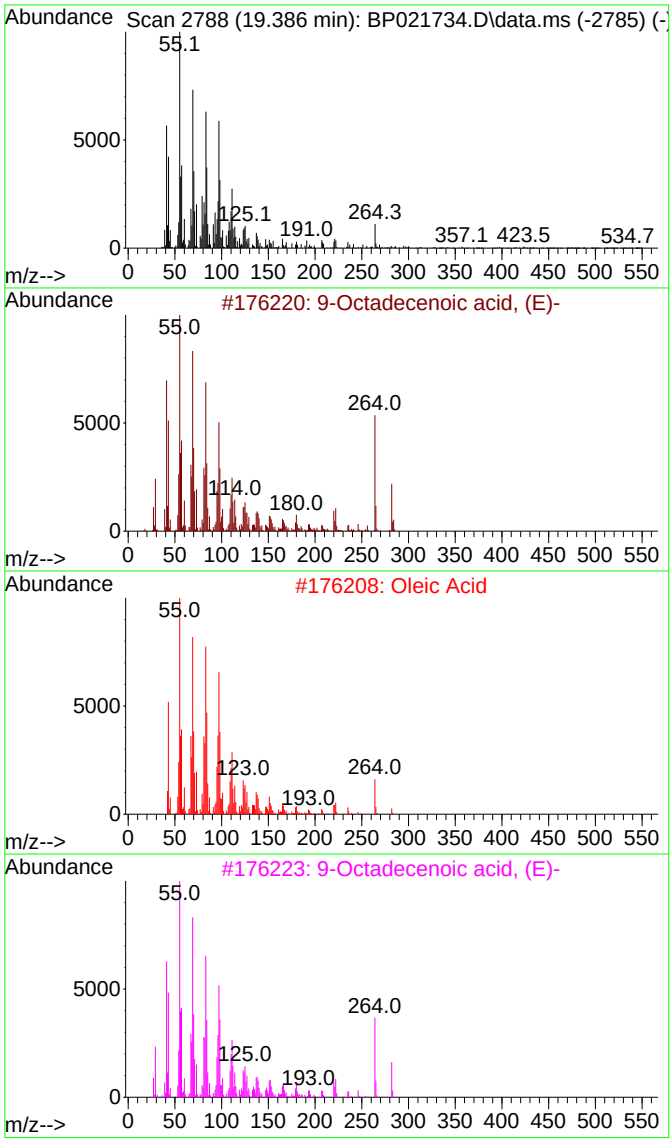
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

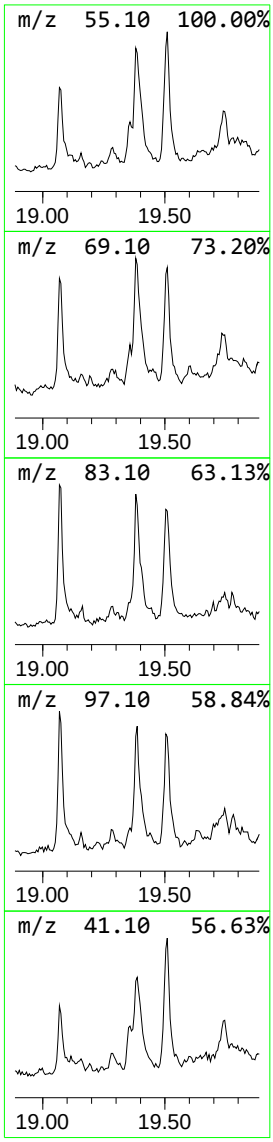
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 9-Octadecenoic acid, (E)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.386	4.52 ng/ul	1502200	Phenanthrene-d10	17.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	97
2	Oleic Acid	282	C18H34O2	000112-80-1	86
3	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	83
4	14-Pentadecenoic acid	240	C15H28O2	017351-34-7	83
5	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	74





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Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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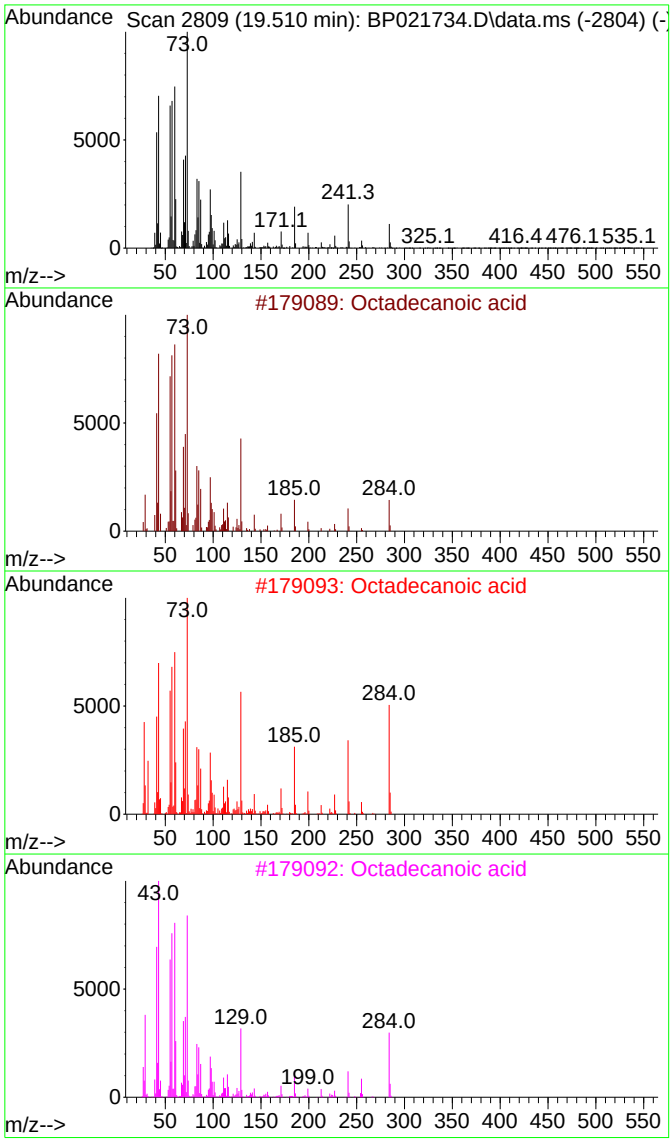
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

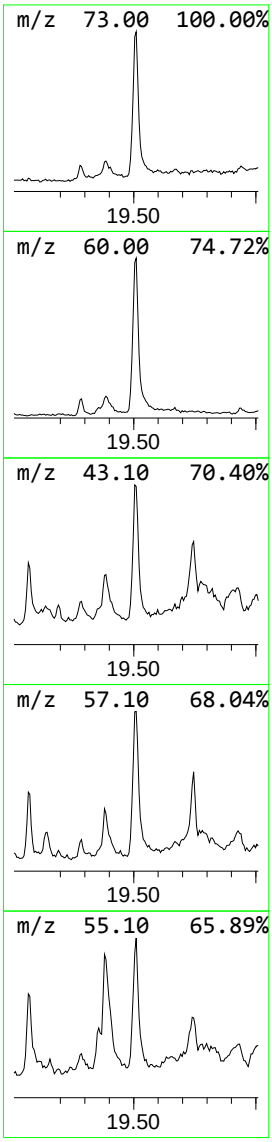
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.510	4.67 ng/ul	1445330	Chrysene-d12	21.686

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	94
4	Octadecanoic acid	284	C18H36O2	000057-11-4	93
5	Octadecanoic acid	284	C18H36O2	000057-11-4	81





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Data File : BP021734.D
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ALS Vial : 22 Sample Multiplier: 1

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BNA_P
ClientSampleId :
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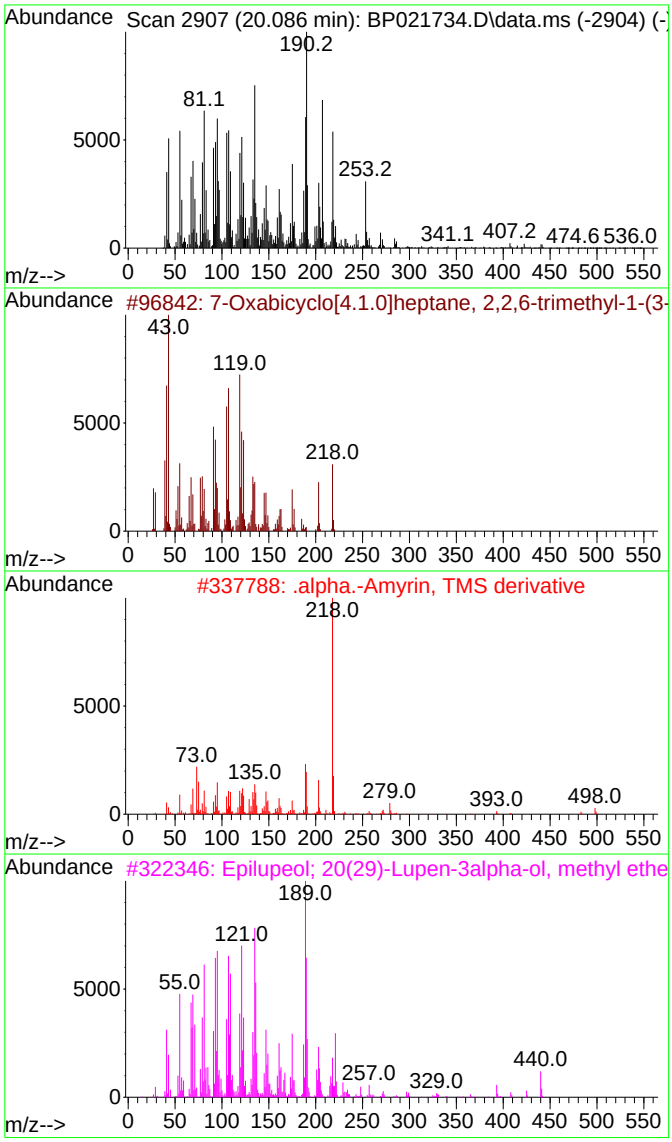
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

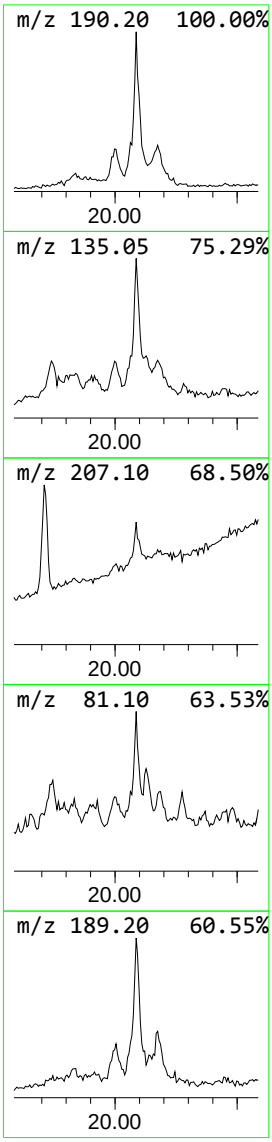
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 7-Oxabicyclo[4.1.0]heptane,... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.086	3.19 ng/ul	987508	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Oxabicyclo[4.1.0]heptane, 2,2,...	218	C15H22O	070038-20-9	62
2	.alpha.-Amyrin, TMS derivative	498	C33H58OSi	1000374-76-6	53
3	Epilupeol; 20(29)-Lupen-3alpha-o...	440	C31H52O	1000513-01-9	45
4	Acetic acid, 2-(2-propyl-1-benzi...	218	C12H14N2O2	331736-92-6	44
5	24-Norursa-3,12-diene	394	C29H46	201358-25-0	41





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Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

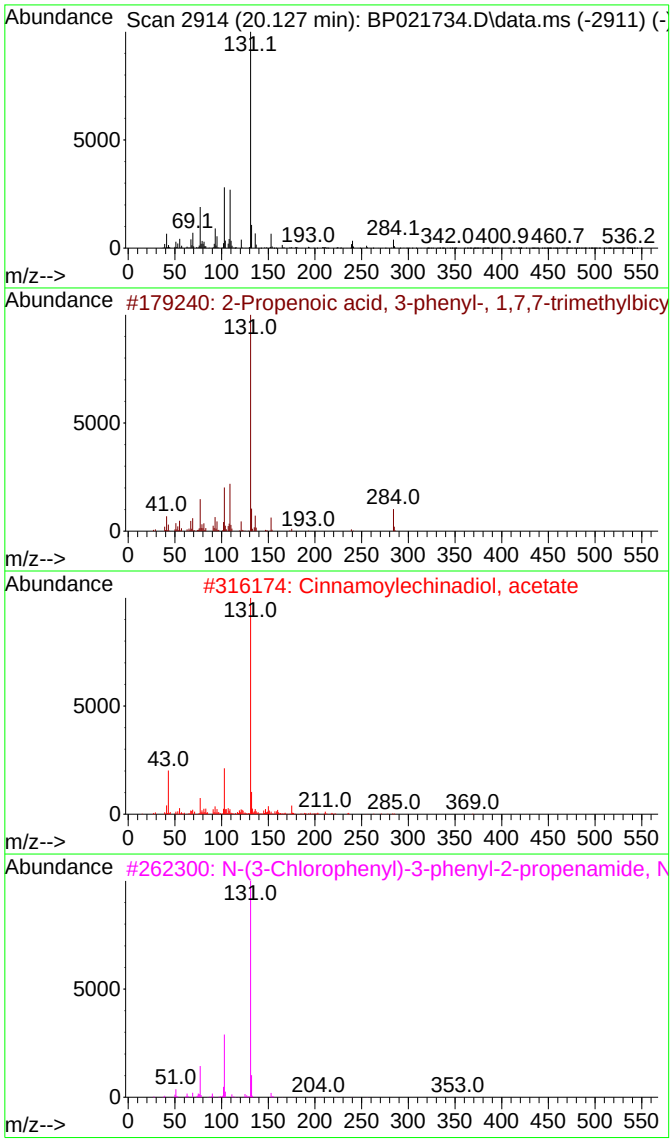
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

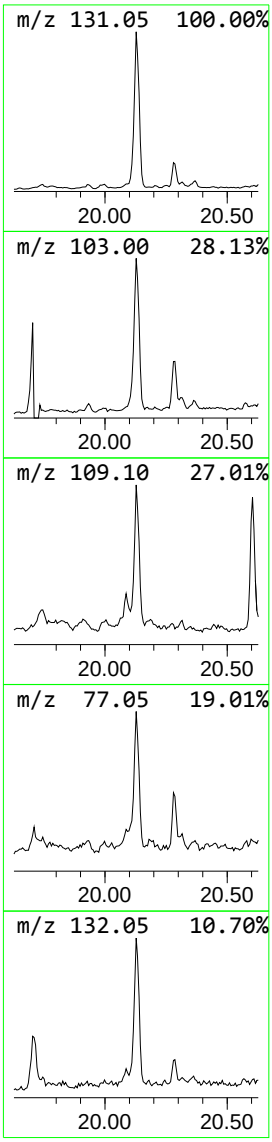
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 2-Propenoic acid, 3-phenyl-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.127	3.64 ng/ul	1125090	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenoic acid, 3-phenyl-, 1,7...	284	C19H24O2	006330-67-2	93
2	Cinnamoylechinadiol, acetate	426	C26H34O5	1000503-42-5	59
3	N-(3-Chlorophenyl)-3-phenyl-2-pr...	353	C17H11ClF3NO2	1000492-37-1	53
4	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	53
5	2-Propenoic acid, 3-phenyl-, 1,7...	284	C19H24O2	006330-67-2	45





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

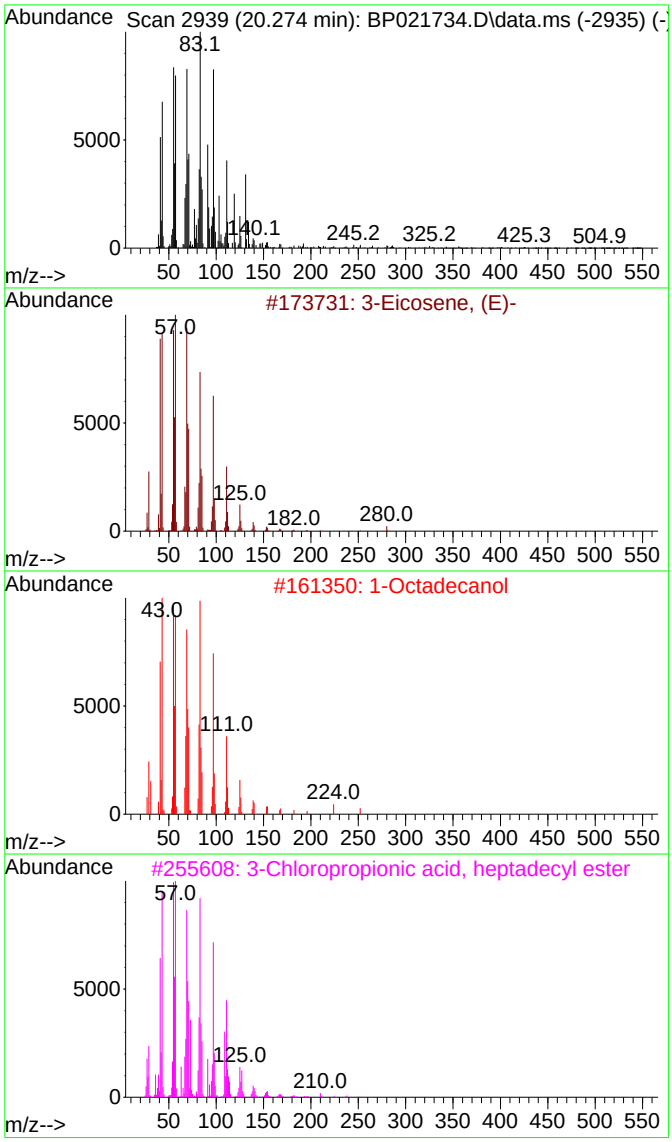
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

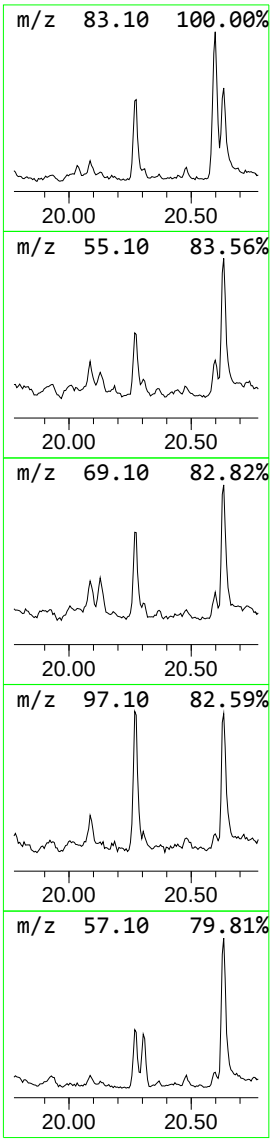
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.275	3.06 ng/ul	946291	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	97
2	1-Octadecanol	270	C18H38O	000112-92-5	90
3	3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	87
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	87
5	1-Heneicosanol	312	C21H44O	015594-90-8	87





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

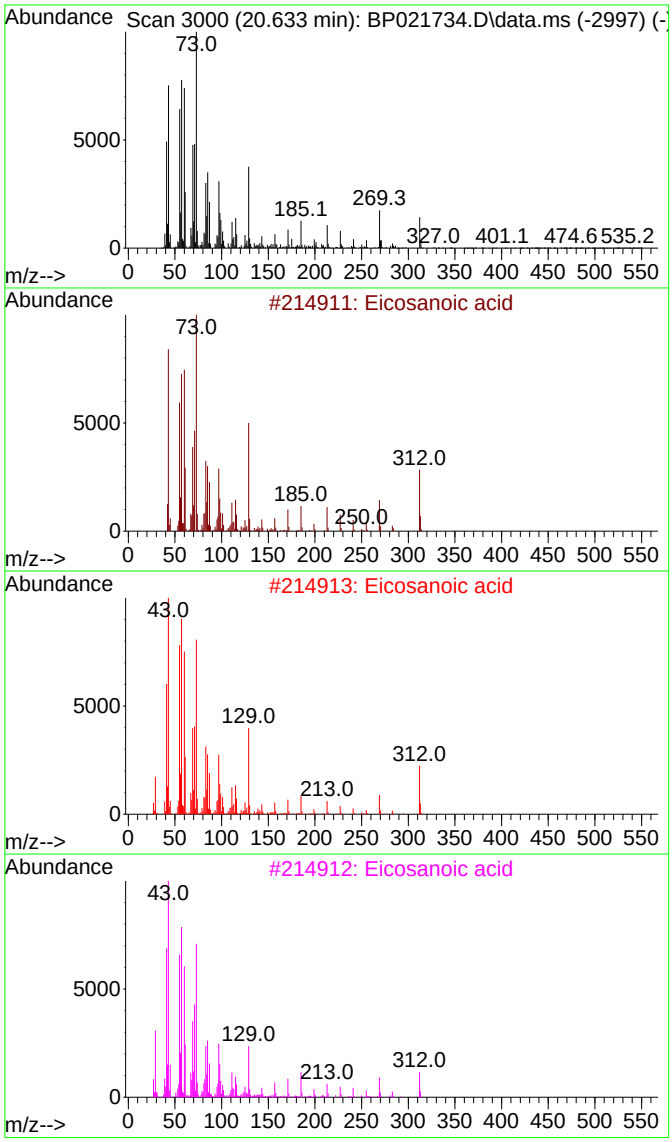
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

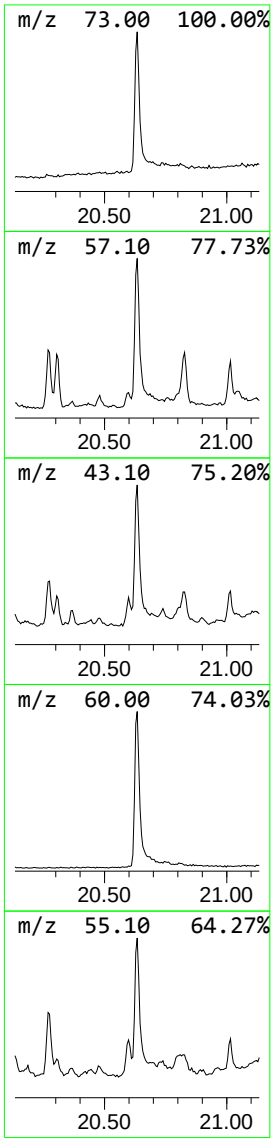
****

Peak Number 18 Eicosanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.633	6.54 ng/ul	2025440	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosanoic acid	312	C20H40O2	000506-30-9	99
2	Eicosanoic acid	312	C20H40O2	000506-30-9	97
3	Eicosanoic acid	312	C20H40O2	000506-30-9	96
4	Heptadecanoic acid	270	C17H34O2	000506-12-7	93
5	Heptadecanoic acid	270	C17H34O2	000506-12-7	91





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

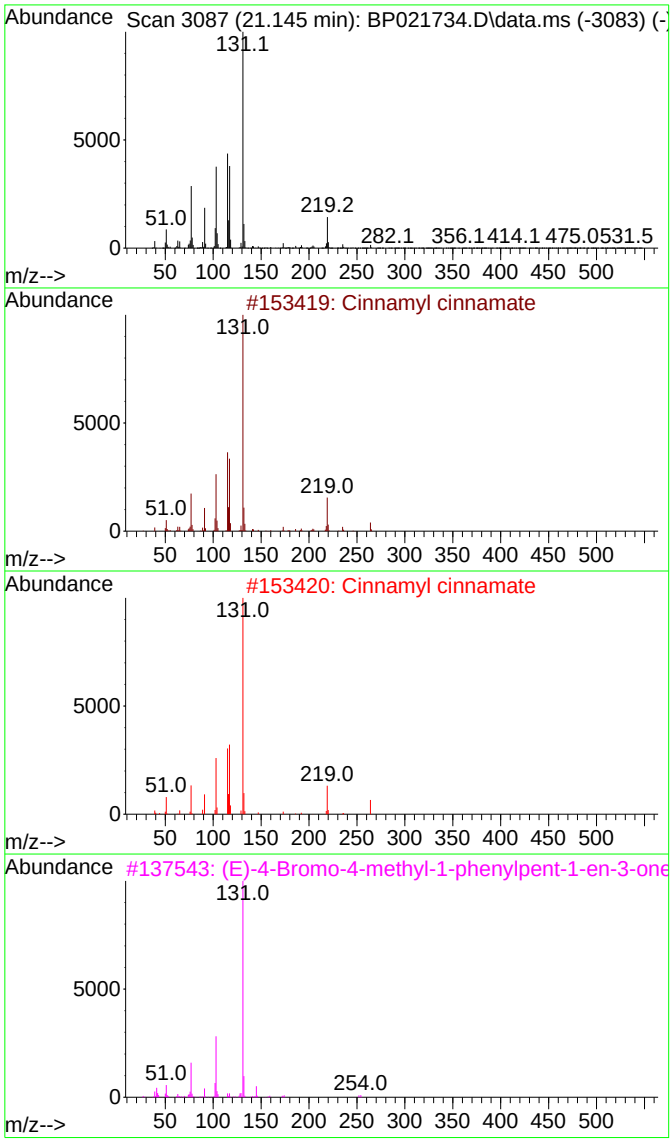
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

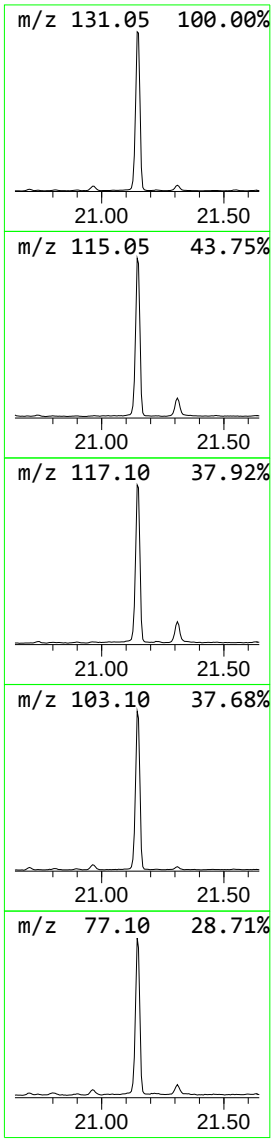
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Cinnamyl cinnamate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.145	31.17 ng/ul	9648140	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	86
3	(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	50
4	(2E)-N-(4-Methylphenyl)-3-phenyl...	237	C16H15NO	134430-88-9	49
5	(2E)-N-(2-Fluorophenyl)-3-phenyl...	241	C15H12FNO	025893-50-9	47





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Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

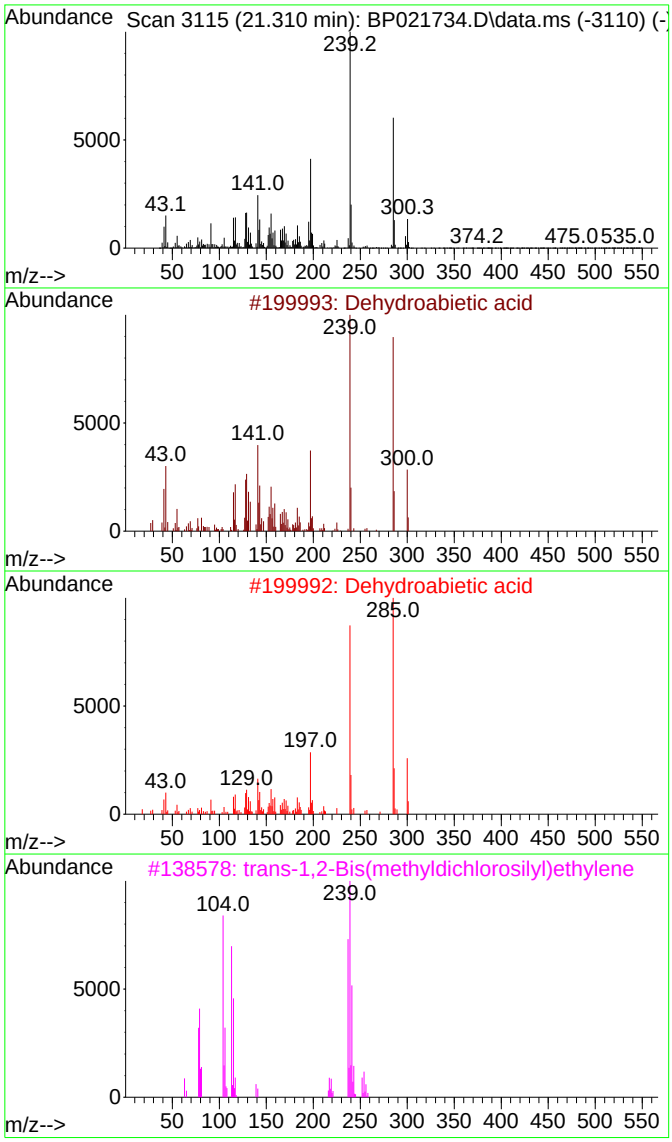
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

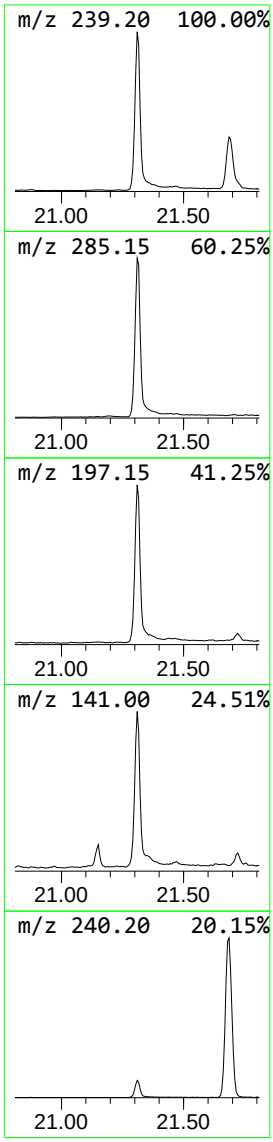
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Dehydroabietic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.310	24.38 ng/ul	7545080	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dehydroabietic acid	300	C20H28O2	001740-19-8	99
2	Dehydroabietic acid	300	C20H28O2	001740-19-8	97
3	trans-1,2-Bis(methyldichlorosily...	252	C4H8Cl4Si2	065899-10-7	94
4	Dehydroabietic acid	300	C20H28O2	001740-19-8	94
5	1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	94





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Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

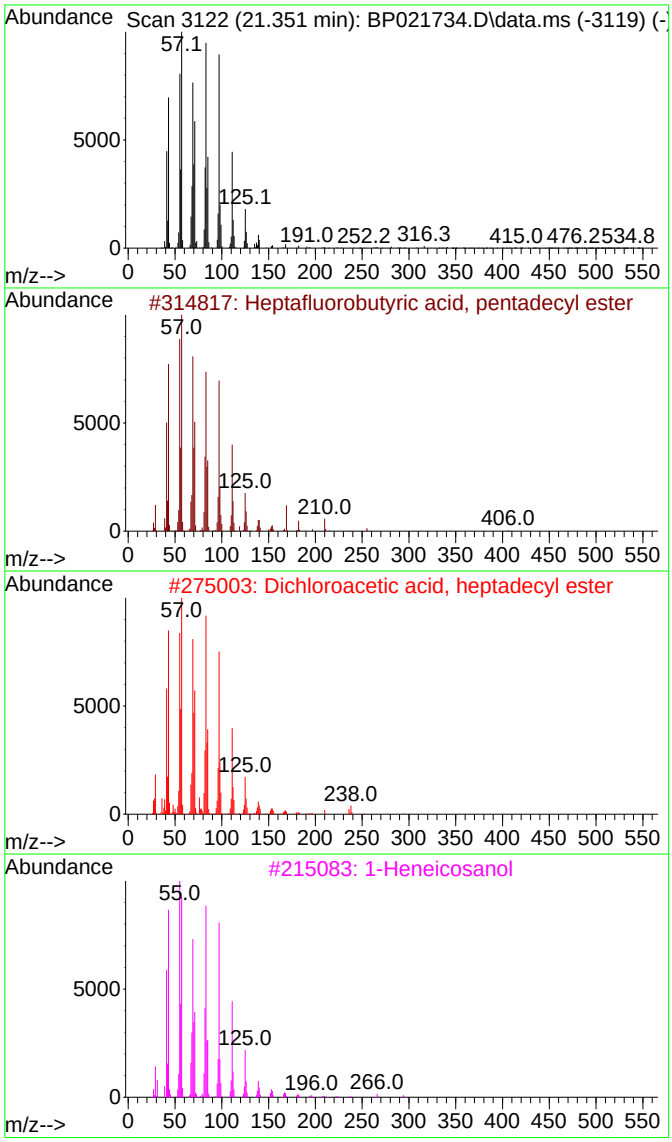
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

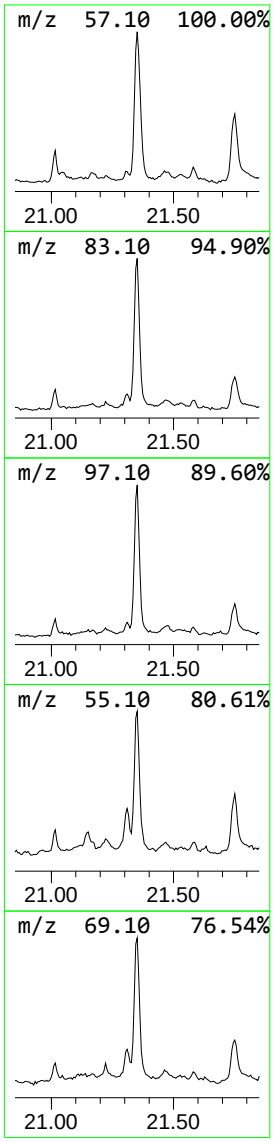
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Heptafluorobutyric acid, pe... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.351	10.39 ng/ul	3215240	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91





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Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

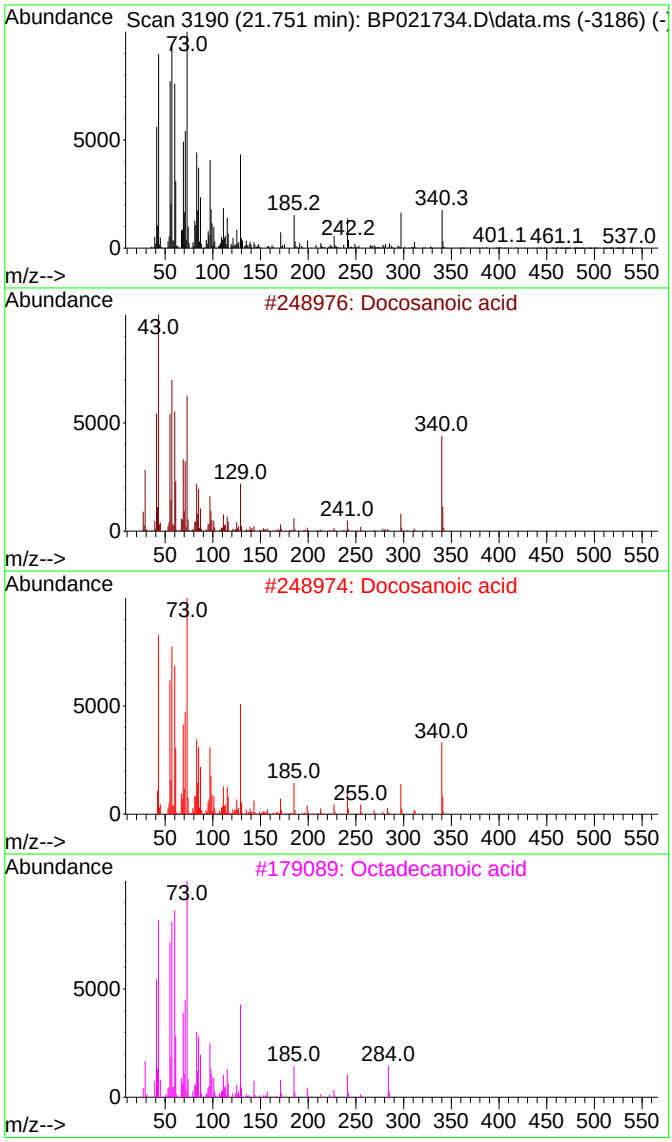
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

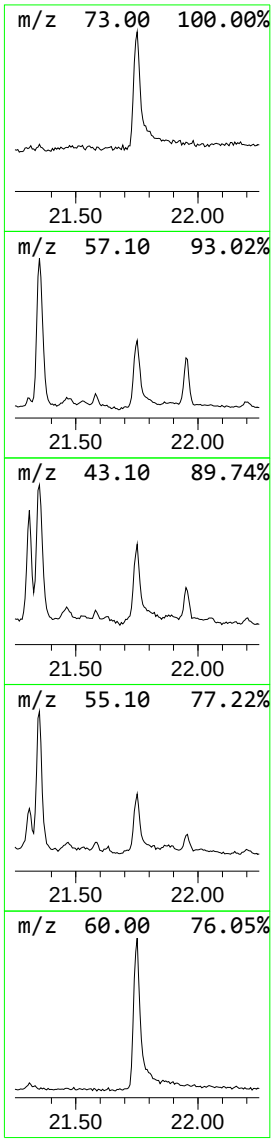
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Docosanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.751	4.59 ng/ul	1420760	Chrysene-d12	21.686

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosanoic acid	340	C22H44O2	000112-85-6	97
2	Docosanoic acid	340	C22H44O2	000112-85-6	96
3	Octadecanoic acid	284	C18H36O2	000057-11-4	72
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	46
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	42





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Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

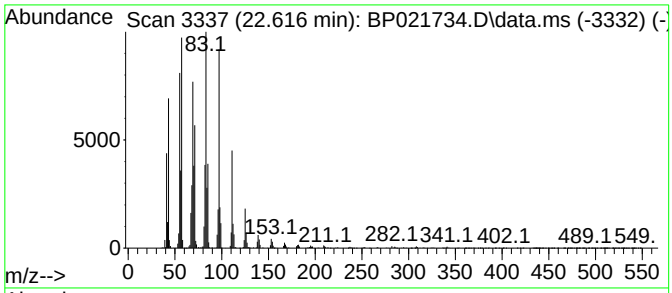
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

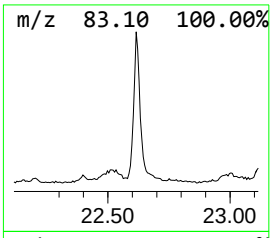
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

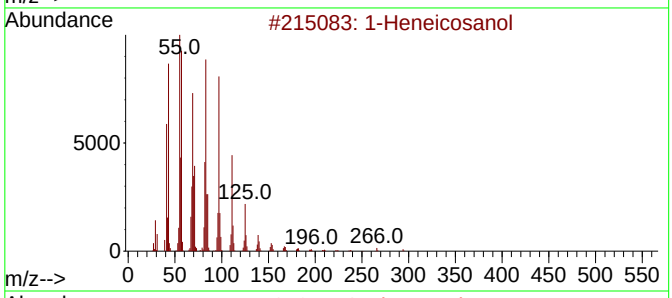
****
Peak Number 23 1-Heneicosanol Concentration Rank 4

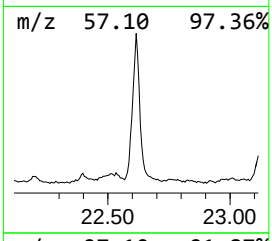
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.616	11.08 ng/ul	3428450	Chrysene-d12	21.686

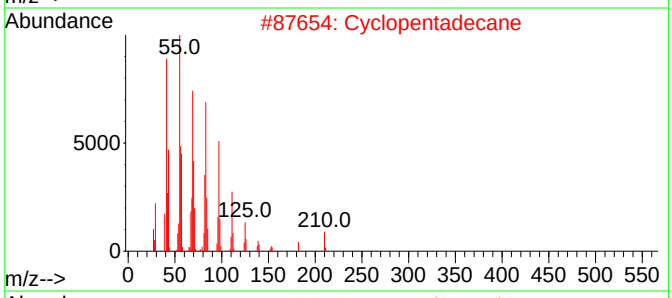
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	93
2	Cyclopentadecane	210	C15H30	000295-48-7	93
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	93
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5	Behenic alcohol	326	C22H46O	000661-19-8	90

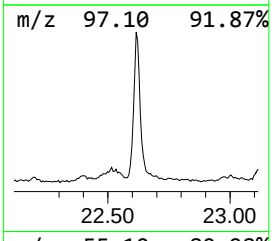


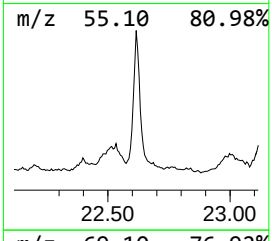


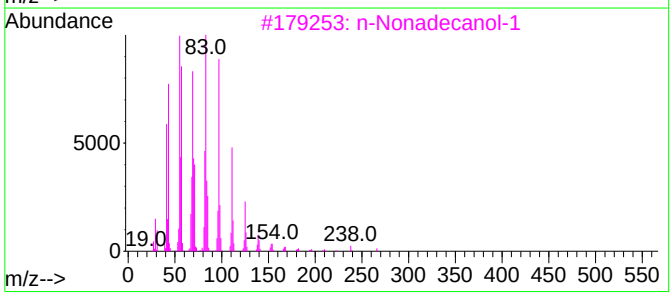


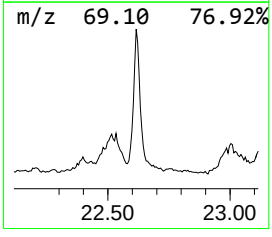












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

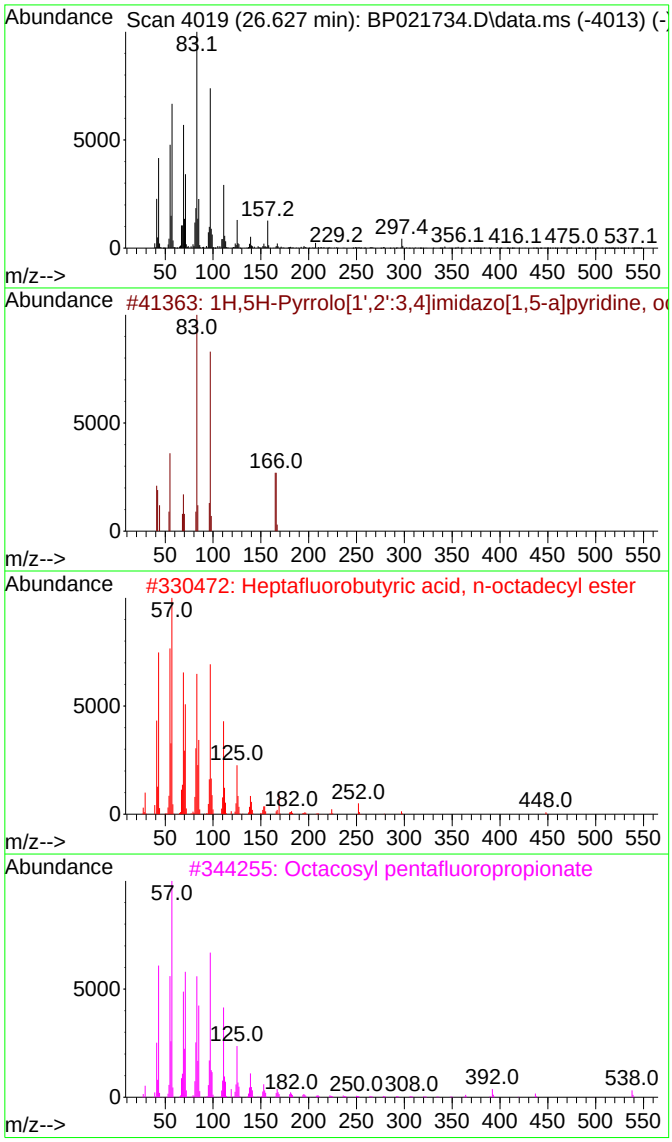
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

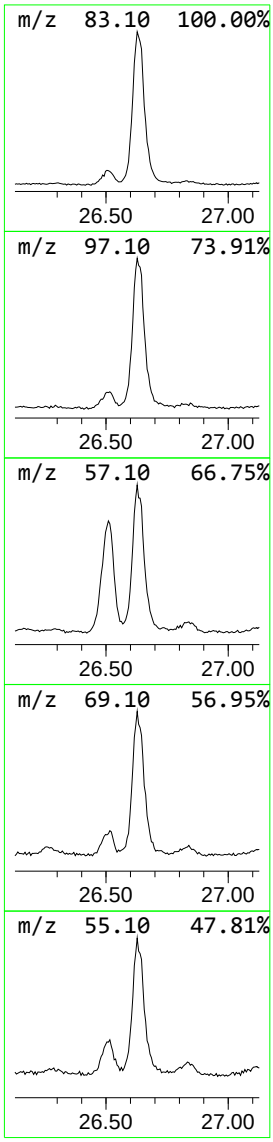
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 1H,5H-Pyrrolo[1',2':3,4]imi... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.627	10.46 ng/ul	3652950	Perylene-d12	25.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H,5H-Pyrrolo[1',2':3,4]imidazo[...	166	C10H18N2	054966-11-9	52
2	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	45
3	Octacosyl pentafluoropropionate	556	C31H57F5O2	1000351-88-9	45
4	Tetracosyl pentafluoropropionate	500	C27H49F5O2	1000351-81-3	43
5	Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	43





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021734.D
Acq On : 30 Aug 2024 00:00
Operator : RC/JU
Sample : P3721-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.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...	3.017	41.3	ng/ul	5707980	1	7.787	2762900	20.0
Propylene Glycol	3.540	2.4	ng/ul	336505	1	7.787	2762900	20.0
.alpha.-Pinene	6.499	3.3	ng/ul	455138	1	7.787	2762900	20.0
Benzoic acid	9.846	4.9	ng/ul	981342	2	10.575	4023780	20.0
1-Phenoxypropan...	11.316	3.5	ng/ul	701890	2	10.575	4023780	20.0
Naphthalene, 1,...	14.522	2.0	ng/ul	535325	3	14.440	5270210	20.0
Naphthalene, 1,...	14.698	2.8	ng/ul	748703	3	14.440	5270210	20.0
.tau.-Cadinol	15.945	2.9	ng/ul	950724	4	17.239	6648690	20.0
(E)-4-(3-Hydrox...	16.616	2.2	ng/ul	717686	4	17.239	6648690	20.0
n-Hexadecanoic ...	18.180	10.0	ng/ul	3329010	4	17.239	6648690	20.0
24-Norursa-3,12...	18.692	2.3	ng/ul	762387	4	17.239	6648690	20.0
(DEL) Alkane: C...	19.075	3.2	ng/ul	1052740	4	17.239	6648690	20.0
9-Octadecenoic ...	19.386	4.5	ng/ul	1502200	4	17.239	6648690	20.0
Octadecanoic acid	19.510	4.7	ng/ul	1445330	5	21.686	6190300	20.0
7-Oxabicyclo[4....	20.086	3.2	ng/ul	987508	5	21.686	6190300	20.0
2-Propenoic aci...	20.127	3.6	ng/ul	1125090	5	21.686	6190300	20.0
(DEL) Alkane: S...	20.275	3.1	ng/ul	946291	5	21.686	6190300	20.0
Eicosanoic acid	20.633	6.5	ng/ul	2025440	5	21.686	6190300	20.0
Cinnamyl cinnamate	21.145	31.2	ng/ul	9648140	5	21.686	6190300	20.0
Dehydroabietic ...	21.310	24.4	ng/ul	7545080	5	21.686	6190300	20.0
Heptafluorobuty...	21.351	10.4	ng/ul	3215240	5	21.686	6190300	20.0
Docosanoic acid	21.751	4.6	ng/ul	1420760	5	21.686	6190300	20.0
1-Heneicosanol	22.616	11.1	ng/ul	3428450	5	21.686	6190300	20.0
1H,5H-Pyrrolo[1...	26.627	10.5	ng/ul	3652950	6	25.086	6982580	20.0