LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCXZ8

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: BP021739.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	28	rVB	19518128	23698156	100.00%	10.478%
2	3.540	91	94	104	rBV	118003	198847	0.84%	0.088%
3	4.922	323	329	335	rBV	106929	163649	0.69%	0.072%
4	5.887	484	493	501	rVB2	70889	140343	0.59%	0.062%
5	6.064	516	523	530	rVB2	75761	127740	0.54%	0.056%
6	6.499	590	597	606	rVB	1406958	2212206	9.33%	0.978%
7	6.805	643	649	651	rBV3	55559	92888	0.39%	0.041%
8	6.840	651	655	662	rVV	119282	217083	0.92%	0.096%
9	6.964	666	676	691	rVB	1196956	2242881	9.46%	0.992%
10	7.122	696	703	714	rVB	957636	1595097	6.73%	0.705%
11	7.246	718	724	731	rBV2	71064	164893	0.70%	0.073%
12	7.328	731	738	745	rVV	1174059	1892732	7.99%	0.837%
13	7.393	745	749	758	rVB3	67354	130356	0.55%	0.058%
14	7.622	781	788	798	rBV5	50795	128807	0.54%	0.057%
15	7.793	809	817	824	rBV	1913023	3221532	13.59%	1.424%
16	8.022	849	856	861	rVV6	49408	120244	0.51%	0.053%
17	8.493	925	936	944	rBV	1290571	2151255	9.08%	0.951%
18	8.605	949	955	964	rVB	511603	889736	3.75%	0.393%
19	8.934	1003	1011	1017	rBV	1104719	1901863	8.03%	0.841%
20	8.975	1017	1018	1024	rVB2	89211	115907	0.49%	0.051%
21	9.493	1099	1106	1113	rBV	121447	216879	0.92%	0.096%
22	9.652	1127	1133	1141	rBV	854145	1457211	6.15%	0.644%
23	9.834	1156	1164	1169	rBV	187917	377223	1.59%	0.167%
24	9.899	1169	1175	1185	rVV5	108227	242678	1.02%	0.107%
25	10.010	1187	1194	1202	rVB	66214	130047	0.55%	0.058%
26	10.199	1218	1226	1246	rBV	1284355	2554925	10.78%	1.130%
27	10.357	1246	1253	1263	rVB3	67091	167223	0.71%	0.074%
28	10.475	1268	1273	1279	rVV2	66128	115338	0.49%	0.051%
29	10.575	1282	1290	1303	rVV	2669974	4699160	19.83%	2.078%
30	10.710	1303	1313	1325	rVB3	93297	284859	1.20%	0.126%
31	11.116	1374	1382	1394	rBV	194621	606731	2.56%	0.268%
32	11.316	1410	1416	1422	rBV	215480	453830	1.92%	0.201%
33	13.169	1725	1731	1736	rBV2	108465	189729	0.80%	0.084%
34	13.334	1755	1759	1765	rVB3	68402	123541	0.52%	0.055%
35	13.698	1816	1821	1830	rBV4	57199	104878	0.44%	0.046%
36	13.834	1837	1844	1852	rBV	2362597	3507129	14.80%	1.551%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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	14.122	1883	1893	1901	rBV	2678997	4283645	18.08%	1.894%
38	14.281	1914	1920	1926	rVB4	62978	125657	0.53%	0.056%
39	14.428	1938	1945	1952	rBV	4211252	6335514	26.73%	2.801%
40	14.510	1955	1959	1966	rVV4	54794	106942	0.45%	0.047%
41	14.640	1974	1981	1998	rBV	1185481	2287967	9.65%	1.012%
42	14.781	2000	2005	2012	rVB	110414	170355	0.72%	0.075%
43	14.857	2012	2018	2022	rBV	146516	234180	0.99%	0.104%
44	15.245	2078	2084	2090	rVB5	46776	94659	0.40%	0.042%
45	15.387	2102	2108	2111	rBV6	54020	112357	0.47%	0.050%
46	15.440	2111	2117	2124	rVB	3529690	5264523	22.21%	2.328%
47	15.551	2129	2136	2144	rBV	776078	1169491	4.93%	0.517%
48	16.339	2265	2270	2275	rBV5	75257	143805	0.61%	0.064%
49	16.622	2311	2318	2323	rBV4	102350	233094	0.98%	0.103%
50	16.875	2357	2361	2369	rVB	145780	260820	1.10%	0.115%
51	17.134	2400	2405	2413	rBV	178786	354672	1.50%	0.157%
52	17.228	2413	2421	2428	rVV	4743361	7624486	32.17%	3.371%
53	17.334	2432	2439	2449	rVB	3859013	5640815	23.80%	2.494%
54	17.416	2450	2453	2459	rVV2	82110	121705	0.51%	0.054%
55	17.534	2469	2473	2483	rBV3	136493	200879	0.85%	0.089%
56	17.892	2530	2534	2538	rBV3	146247	222315	0.94%	0.098%
57	18.045	2554	2560	2566	rVB3	128104	243130	1.03%	0.107%
58	18.104	2567	2570	2575	rBV4	104018	216238	0.91%	0.096%
59	18.163	2575	2580	2593	rVV	2641871	4260800	17.98%	1.884%
60	18.281	2597	2600	2610	rVB2	336477	637273	2.69%	0.282%
61	18.469	2629	2632	2635	rBV	136530	187402	0.79%	0.083%
62	18.516	2635	2640	2645	rVB4	459154	794858	3.35%	0.351%
63	18.633	2654	2660	2663	rBV4	125086	277790	1.17%	0.123%
64	18.675	2664	2667	2674	rVB4	187053	262686	1.11%	0.116%
65	18.975	2713	2718	2724	rBV	244004	444477	1.88%	0.197%
66	19.028	2724	2727	2730	rBV2	90678	119385	0.50%	0.053%
67	19.069	2730	2734	2737	rBV5	117079	224977	0.95%	0.099%
68	19.110	2737	2741	2750	rVV2	878582	1428642	6.03%	0.632%
69	19.281	2767	2770	2774	rVB2	453663	589228	2.49%	0.261%
70	19.375	2777	2786	2797	rBV	2923964	5953909	25.12%	2.633%
71	19.469	2798	2802	2803	rBV	408925	447487	1.89%	0.198%
72	19.498	2803	2807	2818	rVB	1174150	1875526	7.91%	0.829%
73	19.586	2820	2822	2824	rVB	156035	115315	0.49%	0.051%
74	19.622	2824	2828	2836	rBV	766737	1291251	5.45%	0.571%
75	19.698	2836	2841	2847	rBV	3524513	5502121	23.22%	2.433%
76	20.016	2888	2895	2902	rVB4	714099	1443748	6.09%	0.638%

LSC Area Percent Report

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Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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	20.110	2909	2911	2917	rVB	299368	334482	1.41%	0.148%
78	20.169	2918	2921	2925	rBV2	299209	422760	1.78%	0.187%
79	20.292	2939	2942	2947	rBV2	455678	521784	2.20%	0.231%
80	20.510	2975	2979	2986	rBV	1361859	1871740	7.90%	0.828%
81	20.592	2989	2993	2996	rVV2	617249	863077	3.64%	0.382%
82	20.627	2997	2999	3004	rVB2	374049	434016	1.83%	0.192%
83	20.733	3013	3017	3022	rVB	1133704	1660973	7.01%	0.734%
84	20.798	3023	3028	3037	rVV2	1634746	2692885	11.36%	1.191%
85	21.010	3061	3064	3067	rBV2	321050	392084	1.65%	0.173%
86	21.322	3109	3117	3119	rBV	8210511	15087677	63.67%	6.671%
87	21.345	3119	3121	3134	rVB	6284229	9212394	38.87%	4.073%
88	21.627	3165	3169	3173	rBV3	693036	968016	4.08%	0.428%
89	21.680	3173	3178	3186	rVV2	4779369	7602461	32.08%	3.361%
90	21.751	3187	3190	3200	rVB	1430196	2080623	8.78%	0.920%
91	22.204	3264	3267	3275	rVB4	479604	712752	3.01%	0.315%
92	22.627	3332	3339	3354	rBV	9346211	18703426	78.92%	8.270%
93	23.133	3421	3425	3440	rVB2	1152468	2565521	10.83%	1.134%
94	24.251	3608	3615	3620	rBV	1873331	4516930	19.06%	1.997%
95	24.310	3620	3625	3640	rVB	1961901	4810897	20.30%	2.127%
96	24.857	3712	3718	3727	rVB	1272790	3128501	13.20%	1.383%
97	25.098	3751	3759	3769	rVB2	2676398	7602592	32.08%	3.361%
98	26.521	3993	4001	4011	rVB	2051031	5274333	22.26%	2.332%
99	26.645	4013	4022	4040	rVB	2561233	8776079	37.03%	3.880%
100	30.845	4722	4736	4737	rBV2	2943215	7989958	33.72%	3.533%

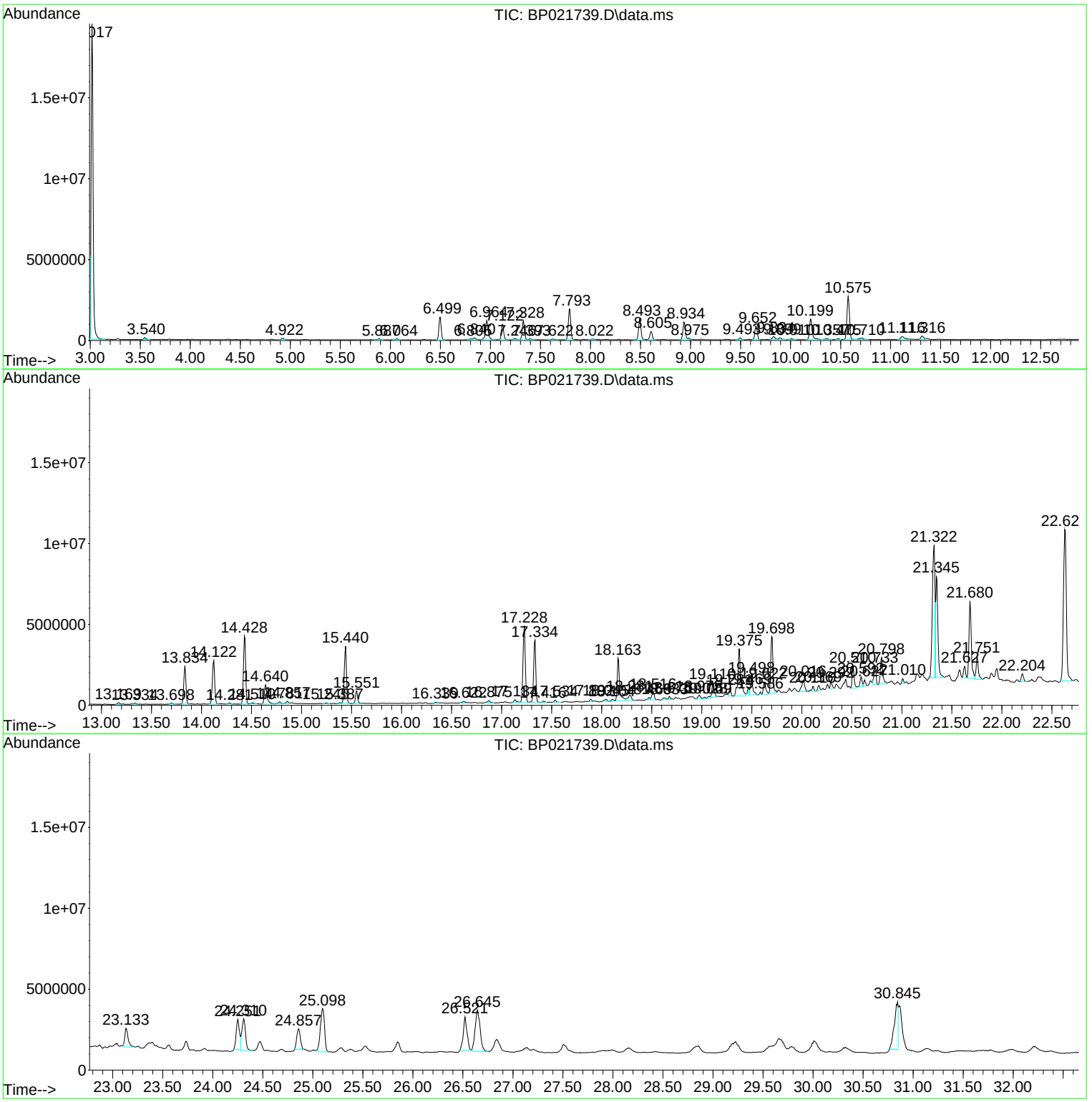
Sum of corrected areas: 226167681

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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 : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

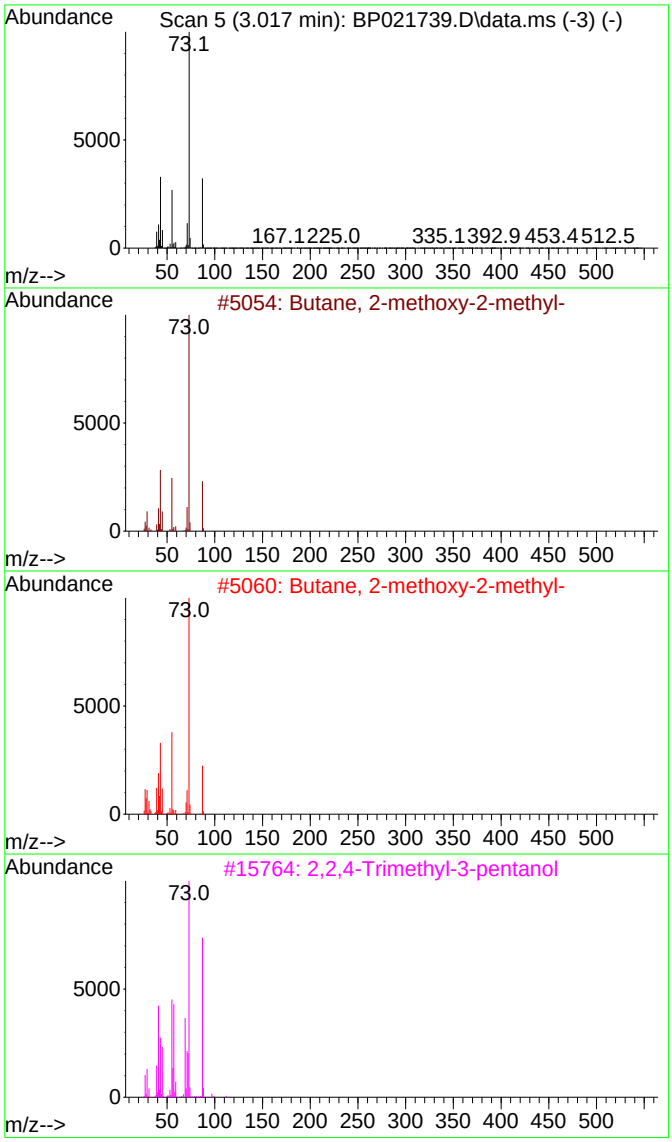
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

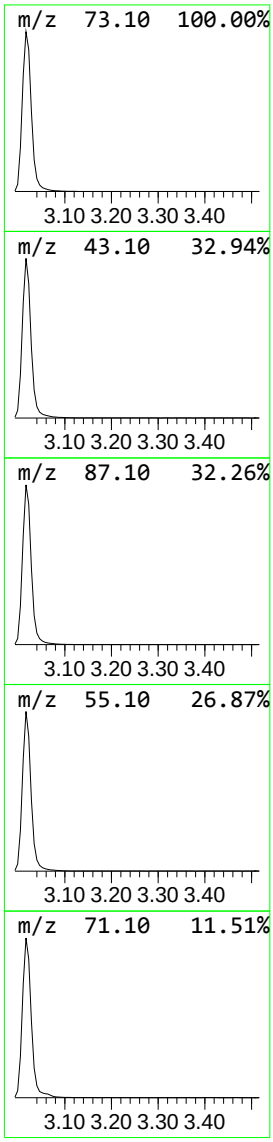
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	147.12 ng/ul	23698200	1,4-Dichlorobenzene-d4	7.793

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	59
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	25
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

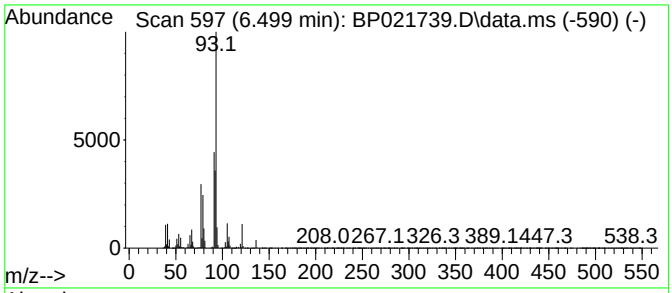
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

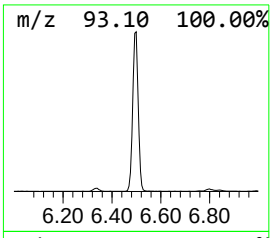
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

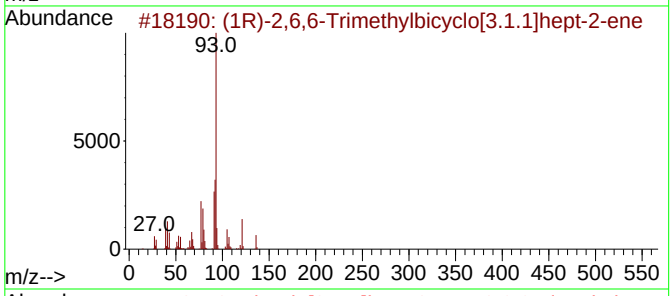
****
Peak Number 2 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 9

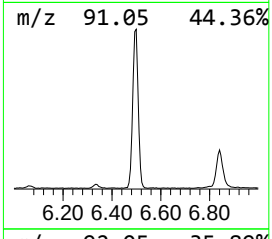
R.T.	EstConc	Area	Relative to ISTD	R.T.
6.499	13.73 ng/ul	2212210	1,4-Dichlorobenzene-d4	7.793

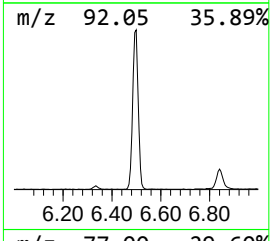
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	94
2	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	94
3	.alpha.-Pinene	136	C10H16	000080-56-8	94
4	.alpha.-Pinene	136	C10H16	000080-56-8	91
5	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91

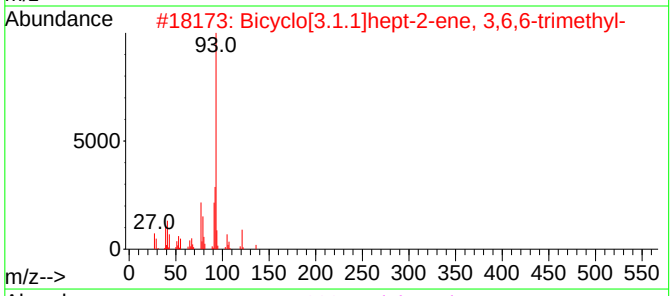


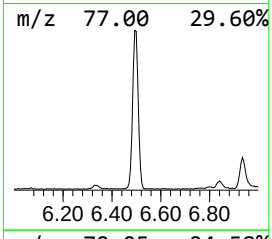


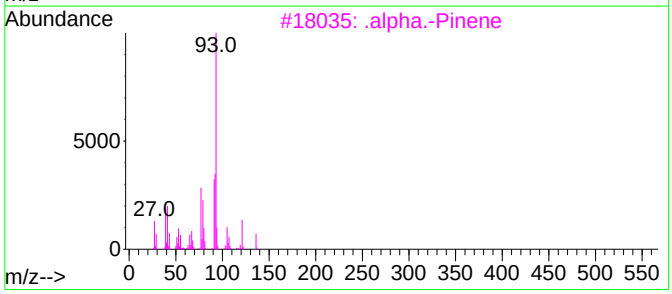


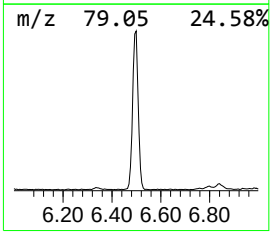












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
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Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

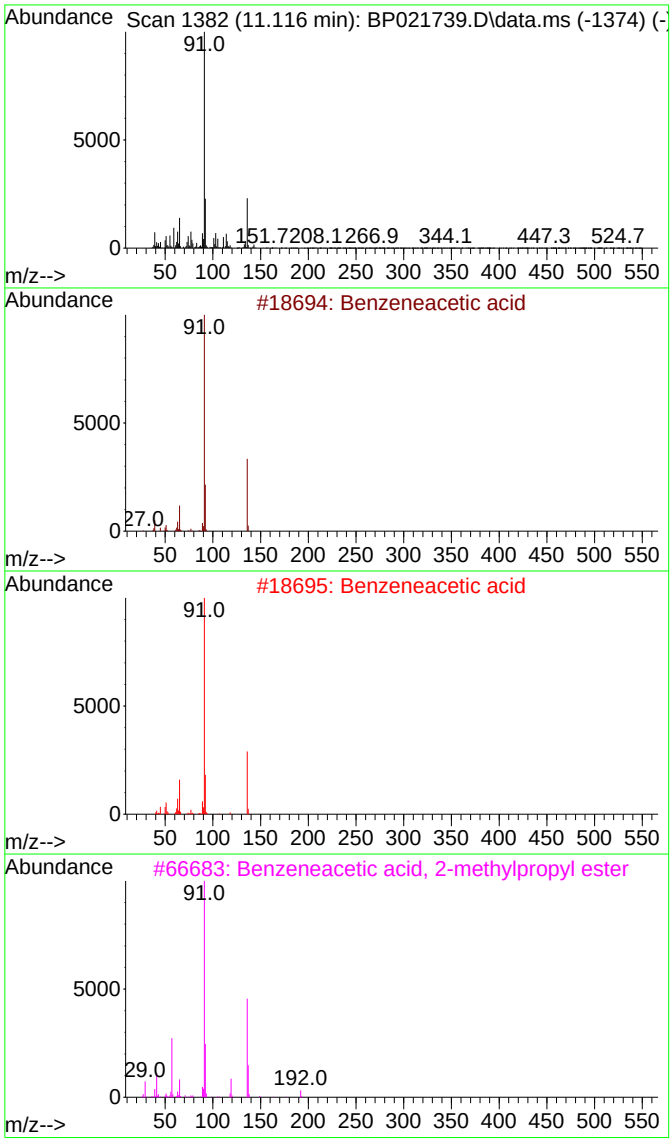
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

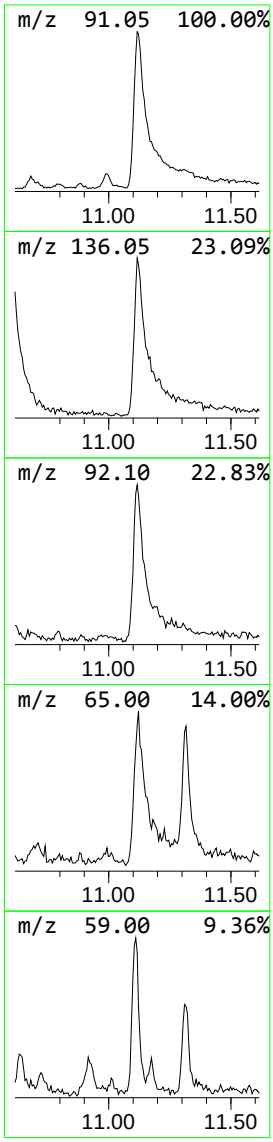
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzeneacetic acid Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.116	2.58 ng/ul	606731	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetic acid	136	C8H8O2	000103-82-2	62
2	Benzeneacetic acid	136	C8H8O2	000103-82-2	53
3	Benzeneacetic acid, 2-methylprop...	192	C12H16O2	000102-13-6	53
4	Benzeneacetic acid	136	C8H8O2	000103-82-2	50
5	Benzeneacetic acid	136	C8H8O2	000103-82-2	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

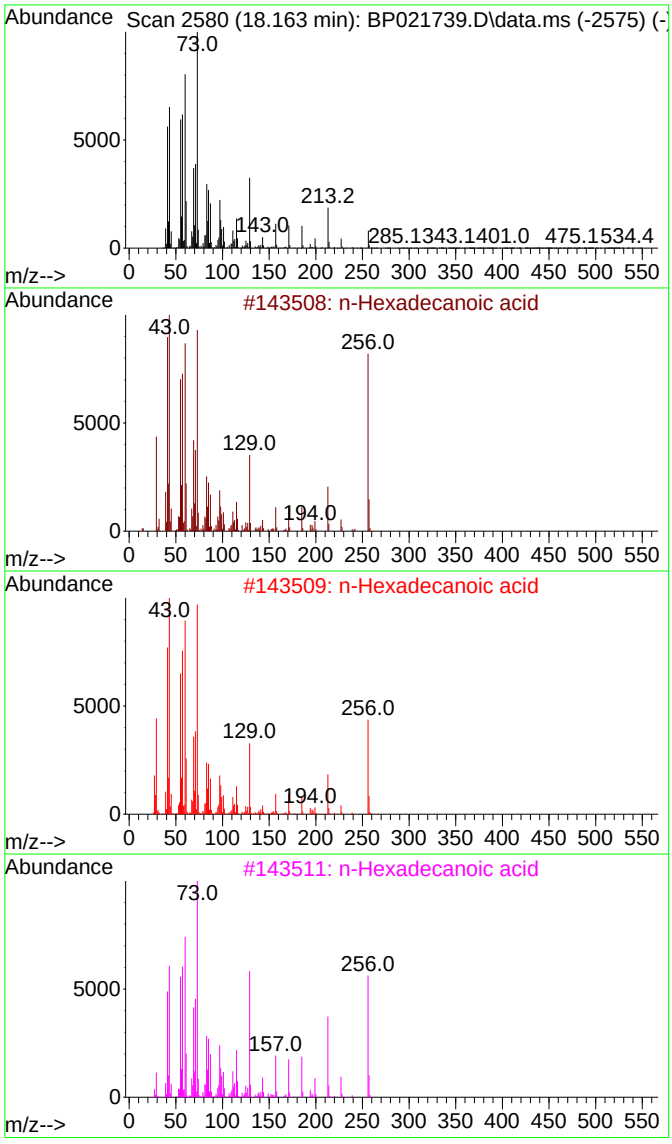
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

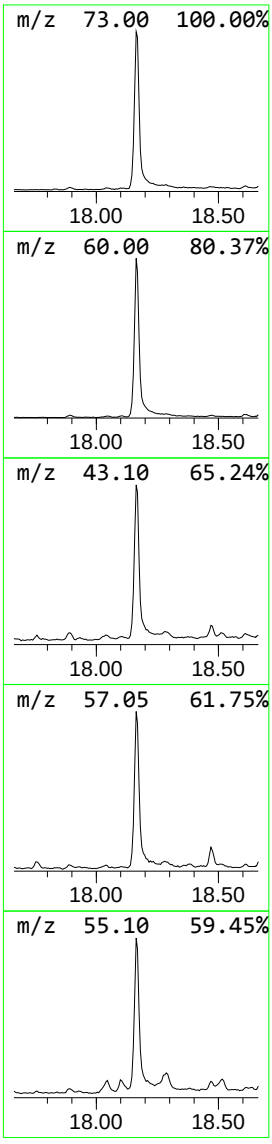
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.163	11.18 ng/ul	4260800	Phenanthrene-d10	17.228

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	98
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 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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

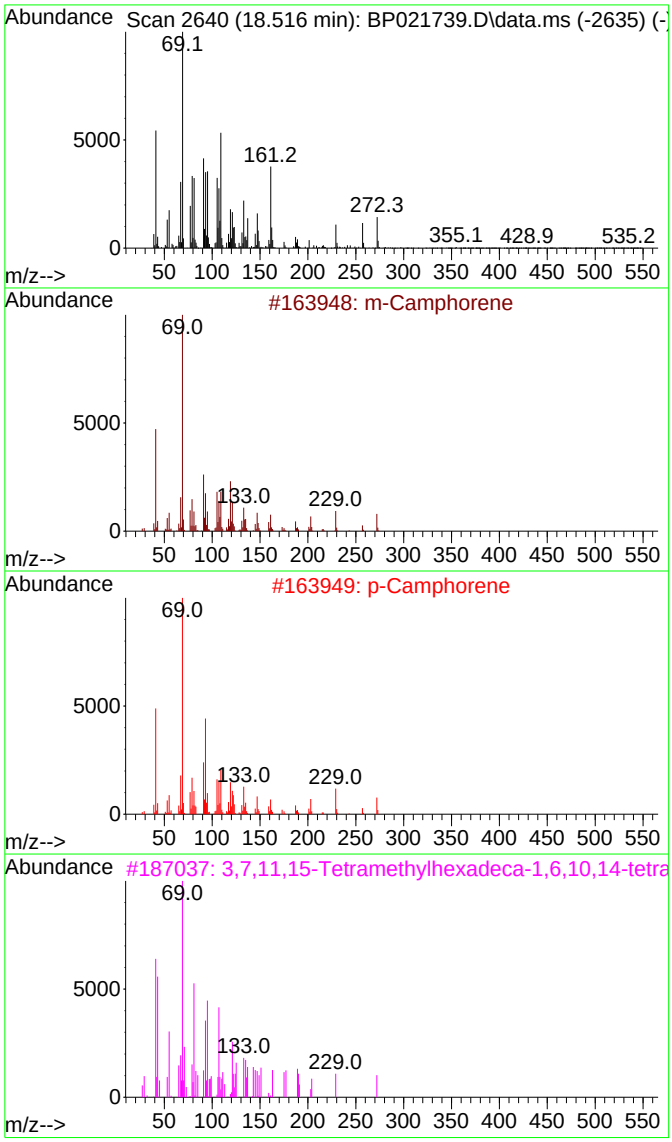
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

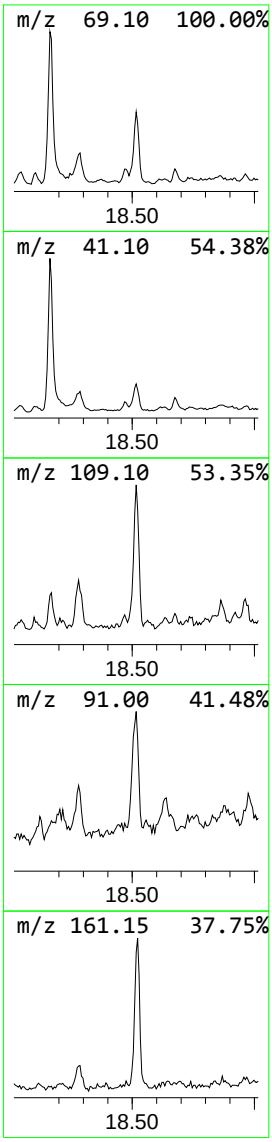
****

Peak Number 5 m-Camphorene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.516	2.09 ng/ul	794858	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m-Camphorene	272	C20H32	020016-73-3	55
2	p-Camphorene	272	C20H32	020016-72-2	53
3	3,7,11,15-Tetramethylhexadeca-1,...	290	C20H34O	068931-30-6	47
4	.beta.-Bisabolenol	220	C15H24O	147126-90-7	43
5	1-Formyl-2,2-dimethyl-3-trans-(3...	220	C15H24O	1000144-09-7	40





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Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCXZ8

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Quant Title : SVOA CALIBRATION

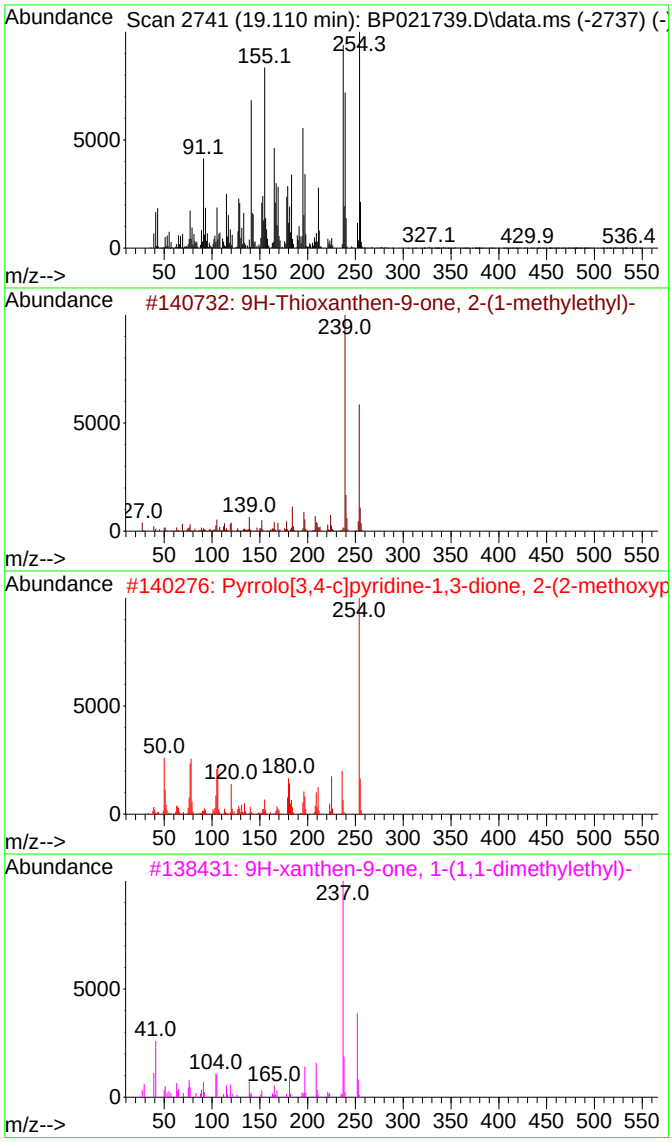
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

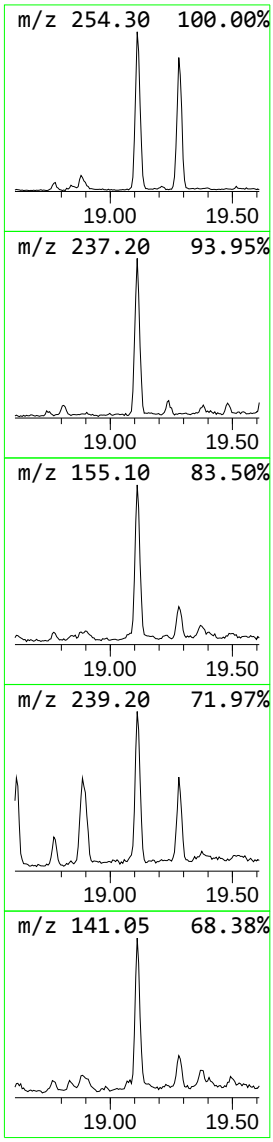
****

Peak Number 6 unknown-01 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.110	3.75 ng/ul	1428640	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Thioxanthen-9-one, 2-(1-methy...	254	C16H14OS	005495-84-1	25
2	Pyrrolo[3,4-c]pyridine-1,3-dione...	254	C14H10N2O3	1000316-41-8	25
3	9H-xanthen-9-one, 1-(1,1-dimethy...	252	C17H16O2	1000396-66-9	25
4	3,6-Dimethoxy-4-phenanthrol	254	C16H14O3	000481-81-2	20
5	Harmol, bis(ethyl)-	254	C16H18N2O	1000395-82-7	20





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Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

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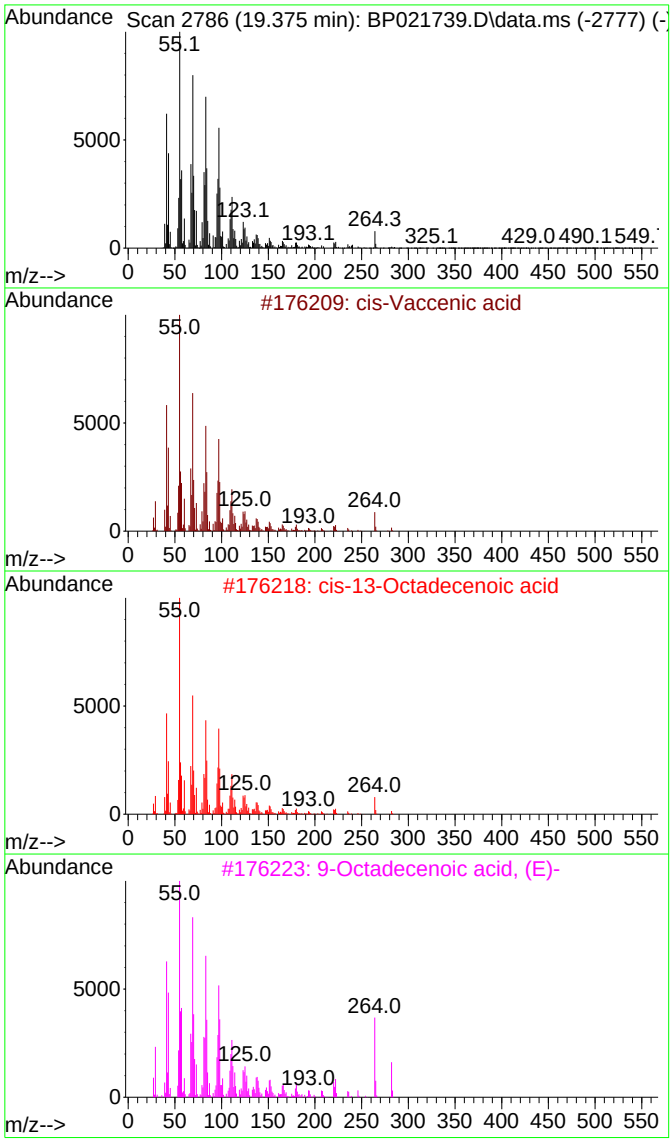
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

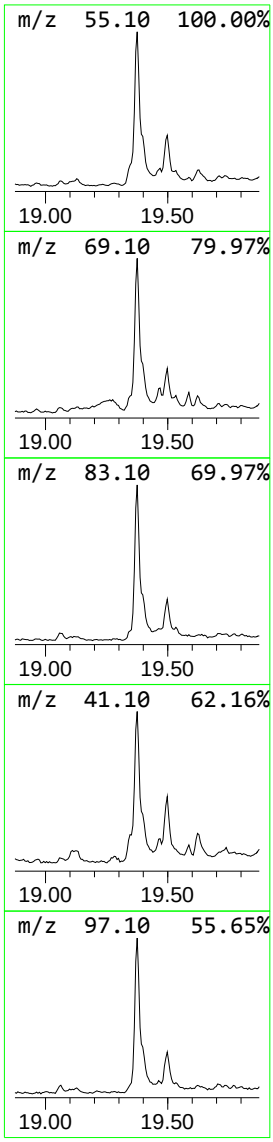
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 cis-Vaccenic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.375	15.62 ng/ul	5953910	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-Vaccenic acid	282	C18H34O2	000506-17-2	97
2	cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	92
3	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	92
4	9-Octadecenoic acid	282	C18H34O2	002027-47-6	91
5	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	91





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Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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

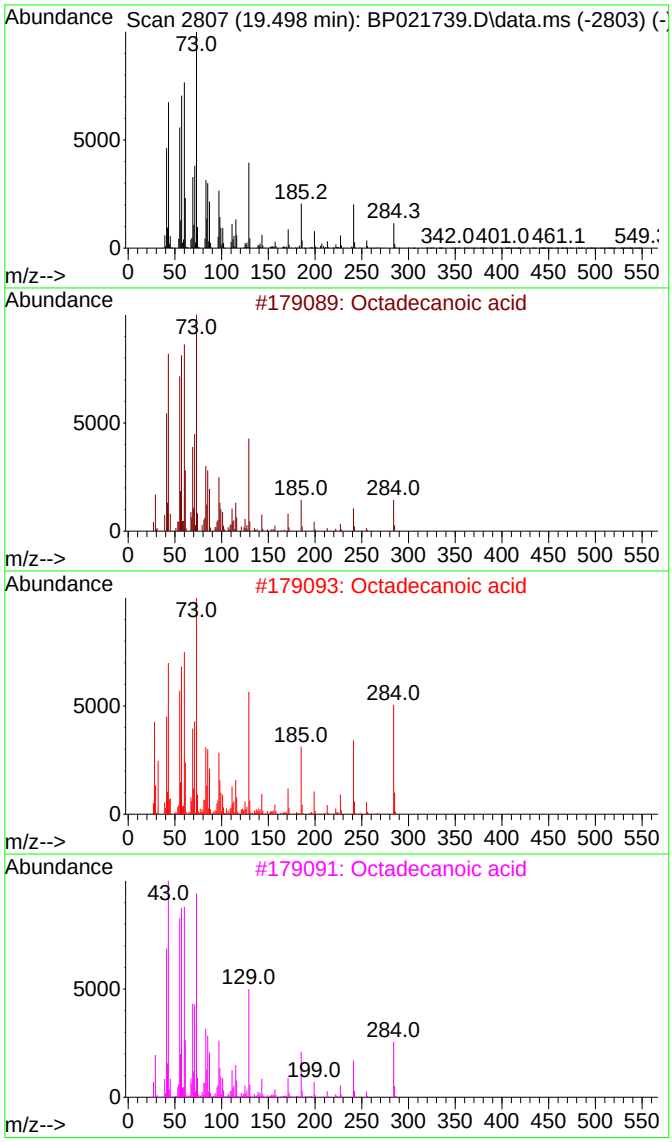
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

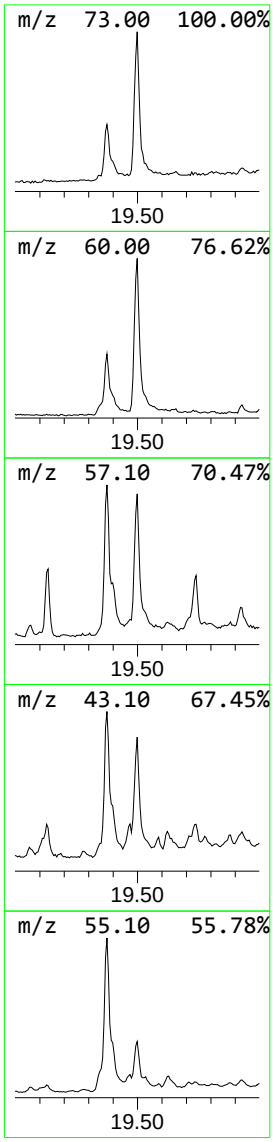
****

Peak Number 8 Octadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.498	4.93 ng/ul	1875530	Chrysene-d12	21.680

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	96
5	Octadecanoic acid	284	C18H36O2	000057-11-4	95





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Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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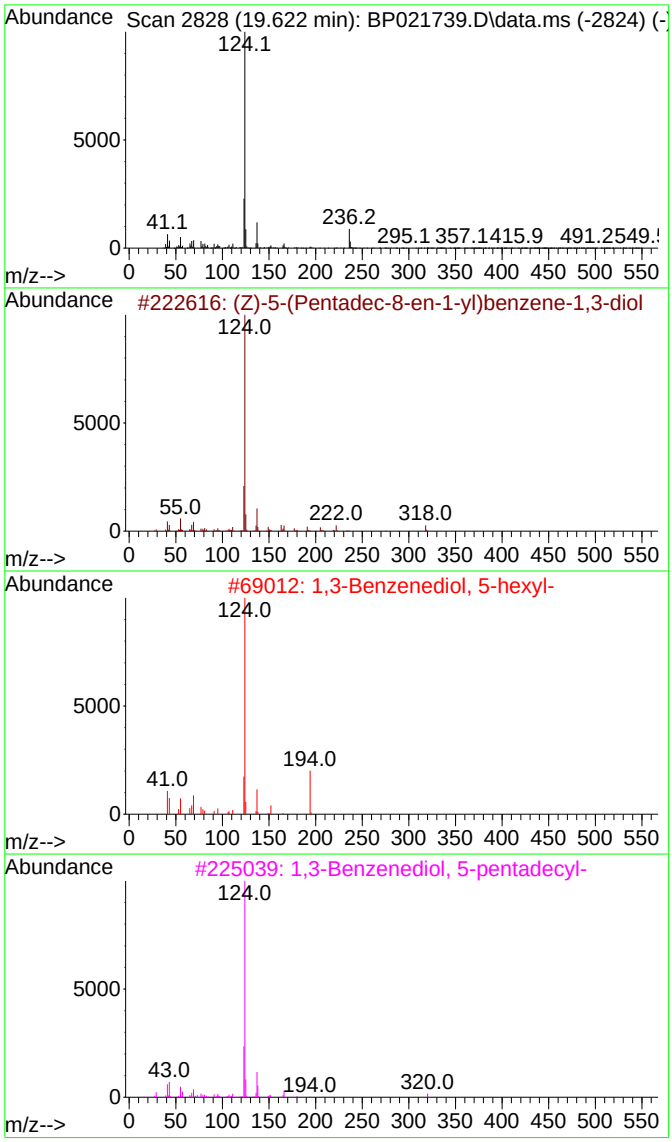
Quant Title : SVOA CALIBRATION

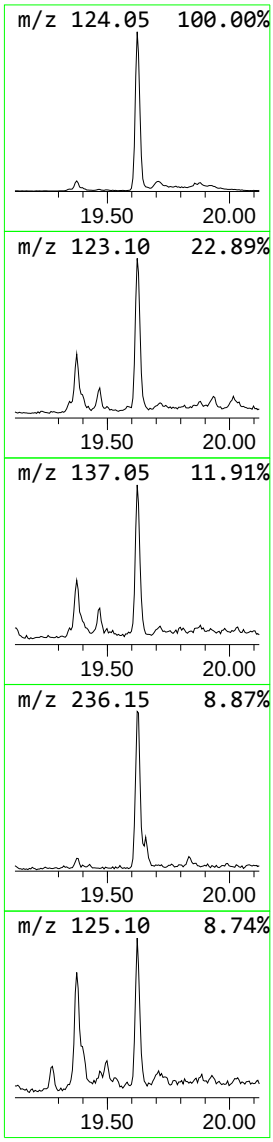
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (Z)-5-(Pentadec-8-en-1-yl)b... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.622	3.40 ng/ul	1291250	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-5-(Pentadec-8-en-1-yl)benzen...	318	C21H34O2	022910-86-7	86
2	1,3-Benzenediol, 5-hexyl-	194	C12H18O2	005465-20-3	83
3	1,3-Benzenediol, 5-pentadecyl-	320	C21H36O2	003158-56-3	80
4	(Z)-5-(Pentadec-8-en-1-yl)benzen...	318	C21H34O2	022910-86-7	78
5	1,3-Benzenediol, 5-pentyl-	180	C11H16O2	000500-66-3	72





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Misc :
ALS Vial : 27 Sample Multiplier: 1

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ClientSampleId :
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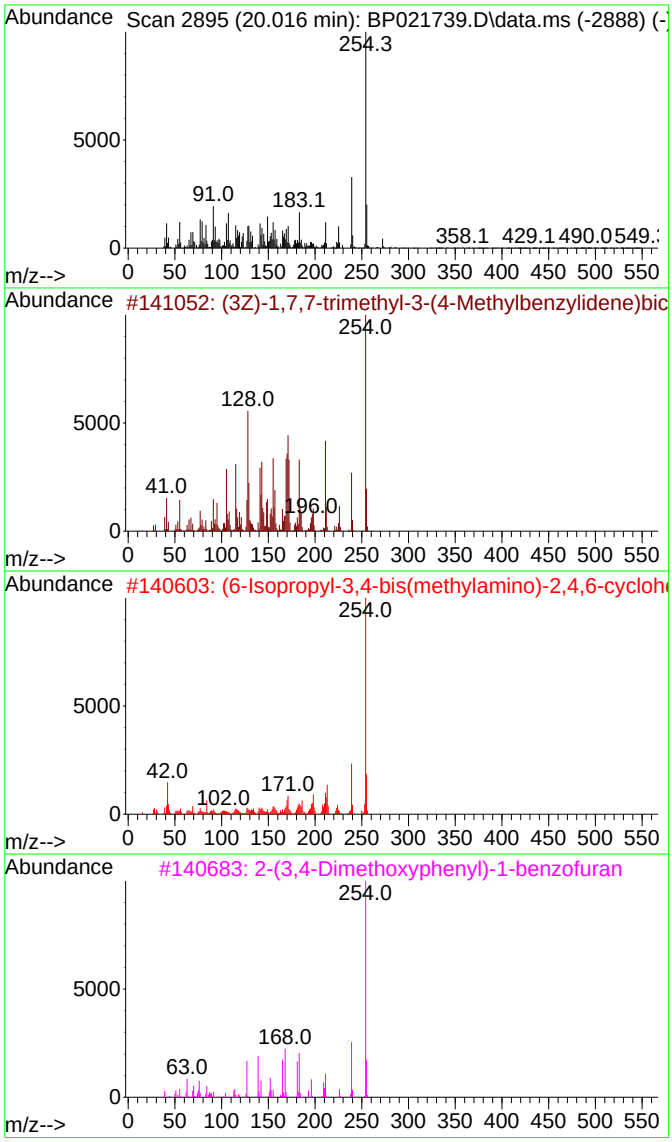
Quant Title : SVOA CALIBRATION

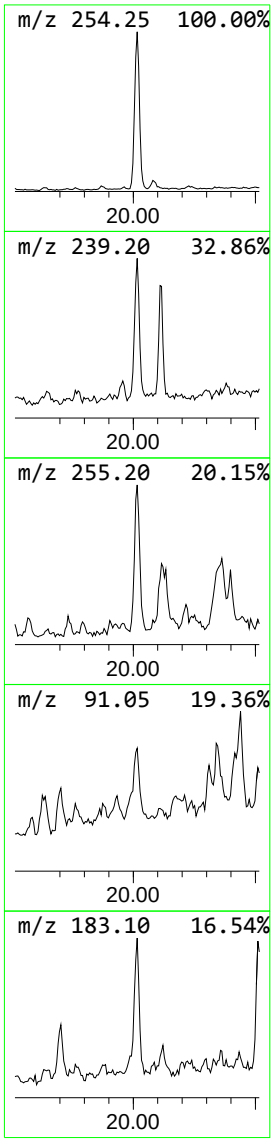
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (3Z)-1,7,7-trimethyl-3-(4-M... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.016	3.80 ng/ul	1443750	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(3Z)-1,7,7-trimethyl-3-(4-Methyl...	254	C18H22O	1000514-88-5	80
2	(6-Isopropyl-3,4-bis(methylamino...	254	C15H18N4	014203-76-0	72
3	2-(3,4-Dimethoxyphenyl)-1-benzof...	254	C16H14O3	051924-87-9	59
4	(E)-4-Hydroxy-4'-methoxychalcone...	296	C18H16O4	1454800-31-7	59
5	1-Propene, 1,2-bis(4-methoxyphen...	254	C17H18O2	020802-02-2	58





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Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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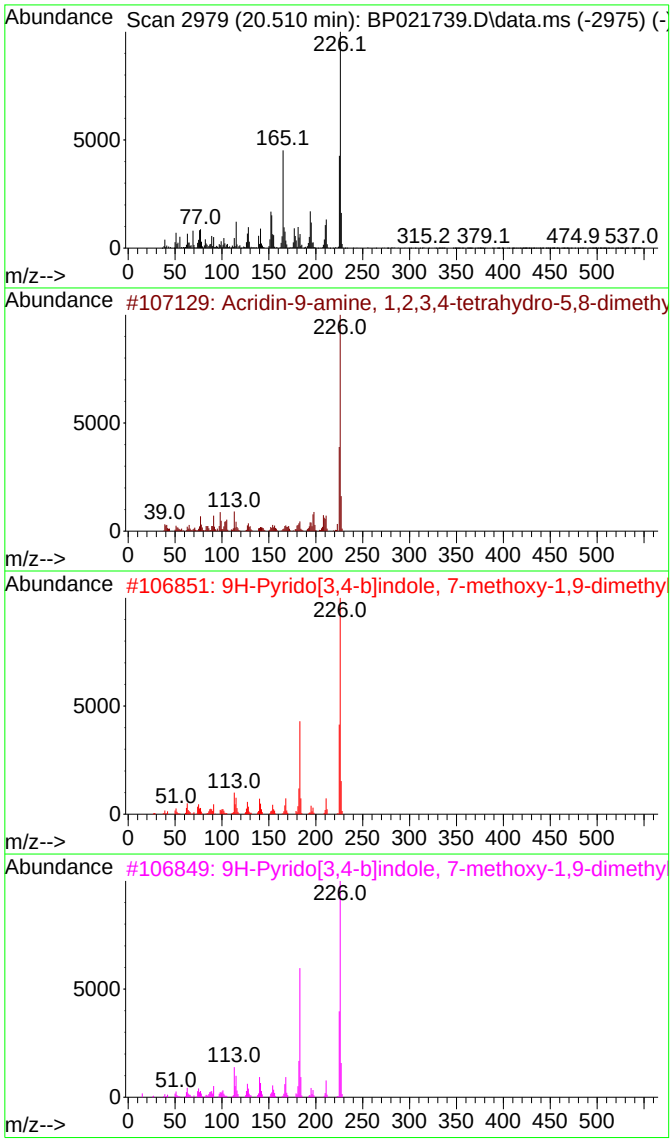
Quant Title : SVOA CALIBRATION

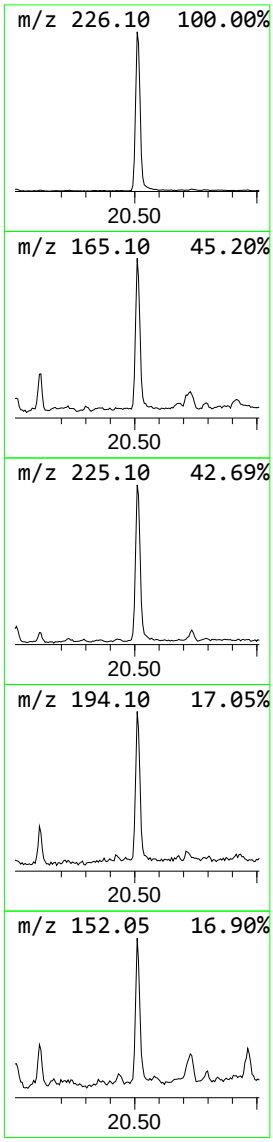
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Acridin-9-amine, 1,2,3,4-te... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.510	4.92 ng/ul	1871740	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	297758-19-1	70
2	9H-Pyrido[3,4-b]indole, 7-methox...	226	C14H14N2O	143502-37-8	52
3	9H-Pyrido[3,4-b]indole, 7-methox...	226	C14H14N2O	143502-37-8	50
4	Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	1000300-57-6	49
5	Benzenamine, N-[(4-nitrophenyl)m...	226	C13H10N2O2	000785-80-8	46





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Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

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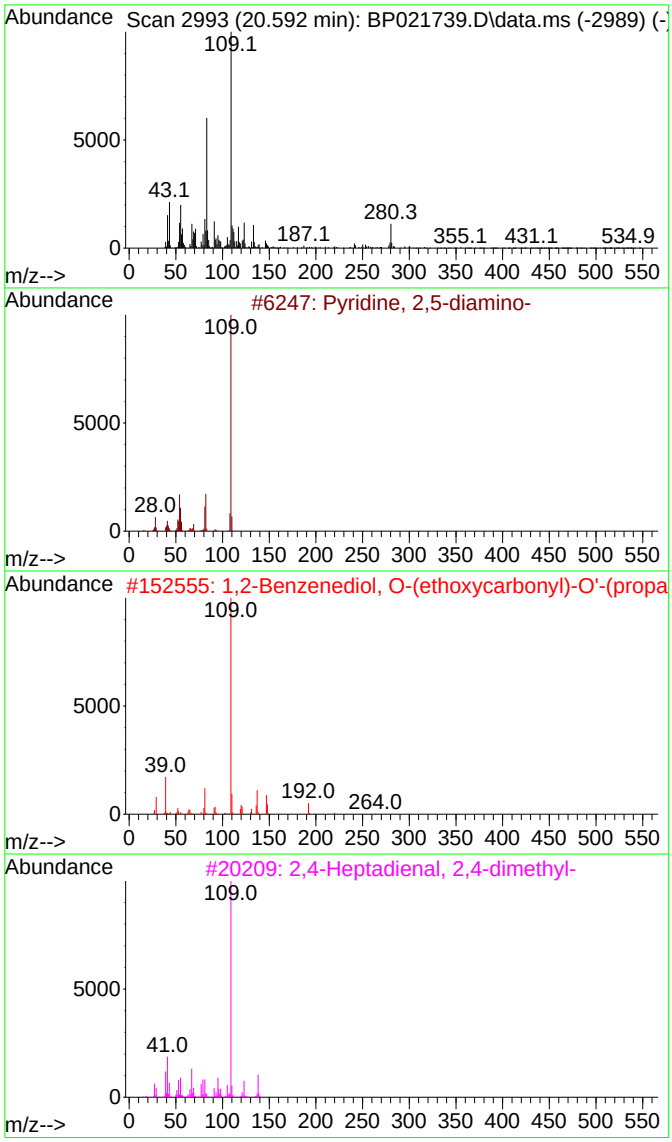
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

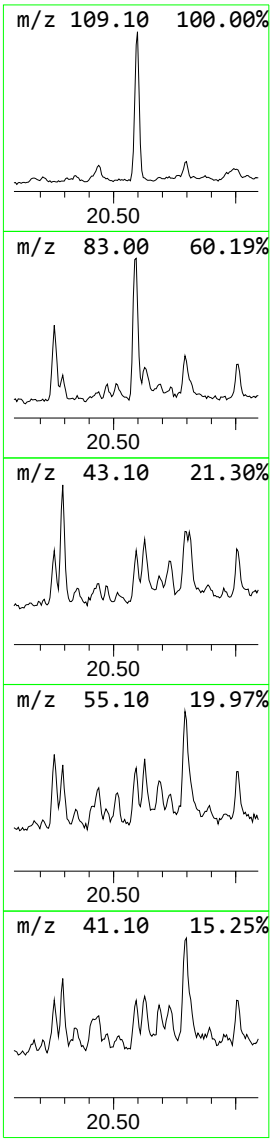
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 unknown-02 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.592	2.27 ng/ul	863077	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2,5-diamino-	109	C5H7N3	004318-76-7	38
2	1,2-Benzenediol, O-(ethoxycarbon...	264	C13H12O6	1000329-71-6	38
3	2,4-Heptadienal, 2,4-dimethyl-	138	C9H14O	042452-48-2	35
4	2-Cyclopentene-1-carboxylic acid...	182	C11H18O2	029915-95-5	35
5	1-Allyl-cyclopropanecarboxylic a...	328	C22H32O2	108546-69-6	35





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Misc :
ALS Vial : 27 Sample Multiplier: 1

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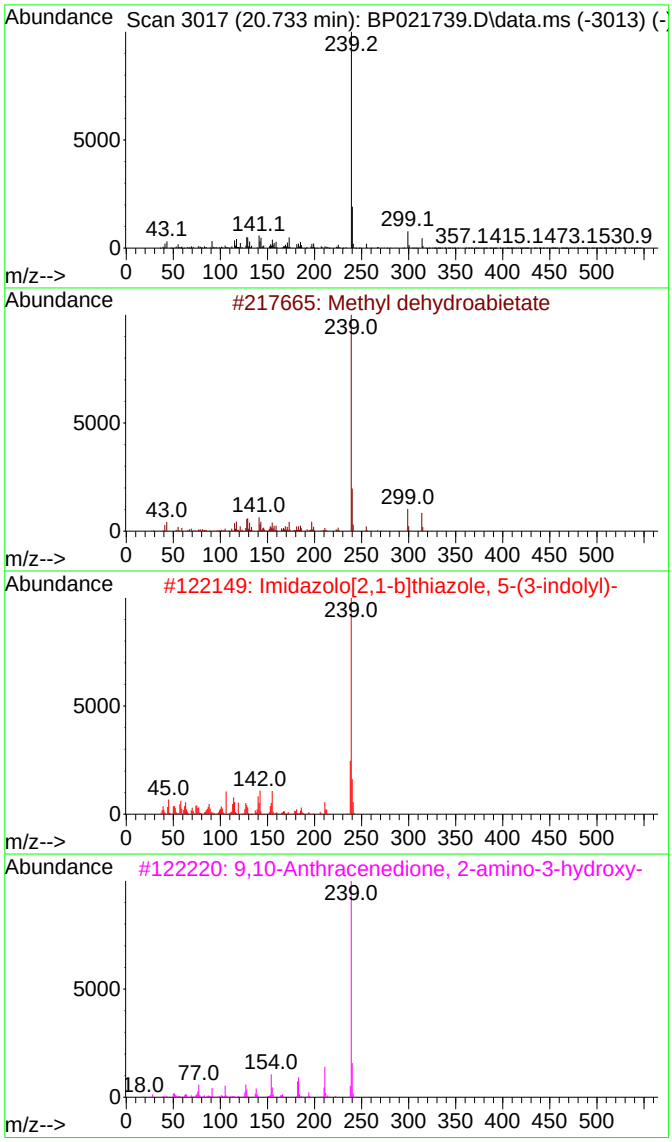
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

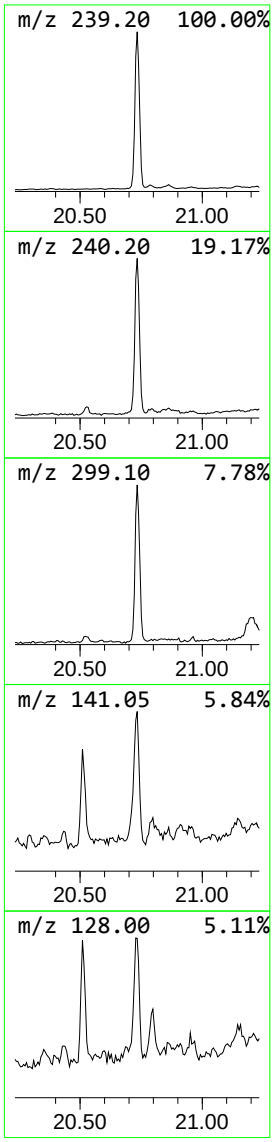
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Methyl dehydroabietate Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.733	4.37 ng/ul	1660970	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl dehydroabietate	314	C21H30O2	001235-74-1	96
2	Imidazolo[2,1-b]thiazole, 5-(3-i...	239	C13H9N3S	327097-37-0	64
3	9,10-Anthracenedione, 2-amino-3-...	239	C14H9NO3	000117-77-1	64
4	9,10-Anthracenedione, 1-amino-4-...	239	C14H9NO3	000116-85-8	59
5	Methyl dehydroabietate	314	C21H30O2	001235-74-1	55





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ALS Vial : 27 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

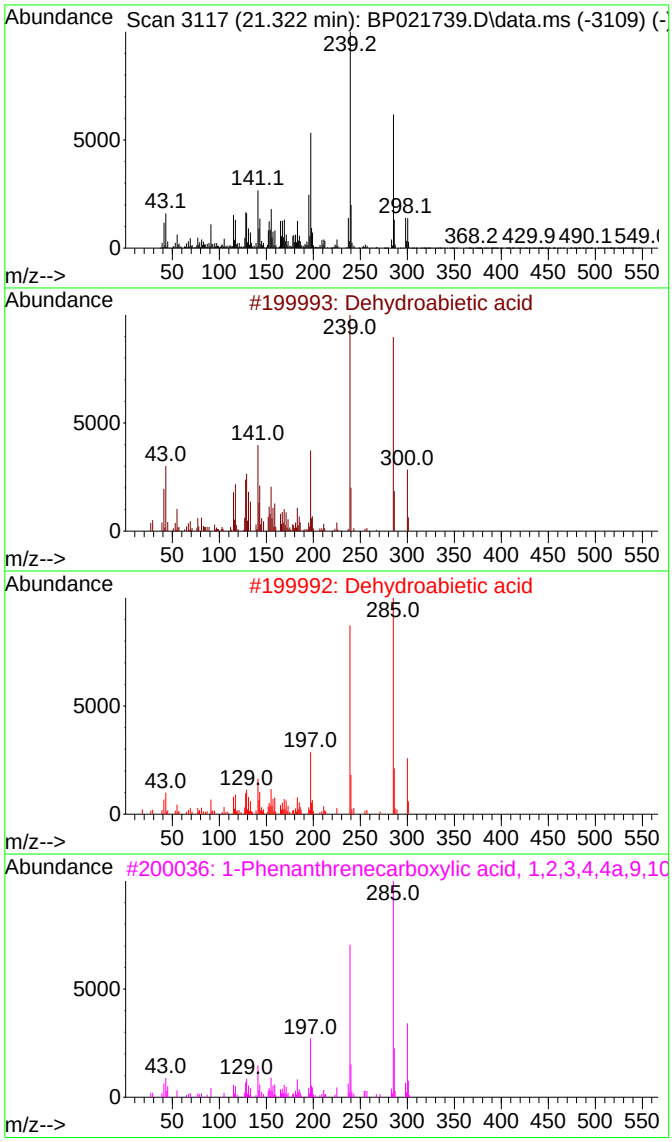
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

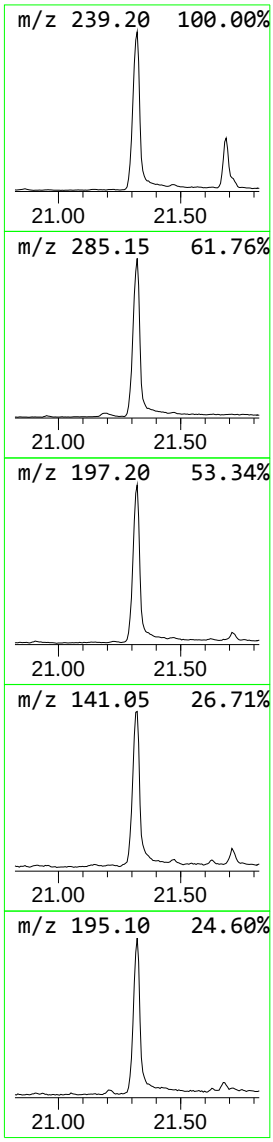
****

Peak Number 15 Dehydroabietic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.322	39.69 ng/ul	15087700	Chrysene-d12	21.680

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	96
3	1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	94
4	trans-1,2-Bis(methyldichlorosily...	252	C4H8Cl4Si2	065899-10-7	93
5	Dehydroabietic acid	300	C20H28O2	001740-19-8	89





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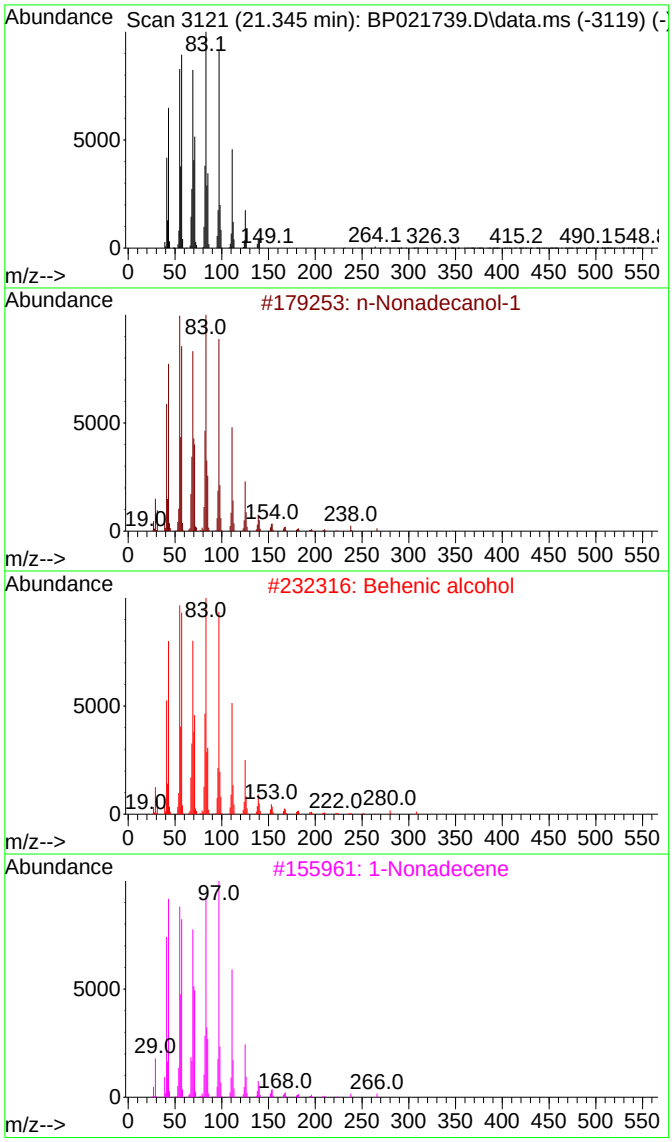
Quant Title : SVOA CALIBRATION

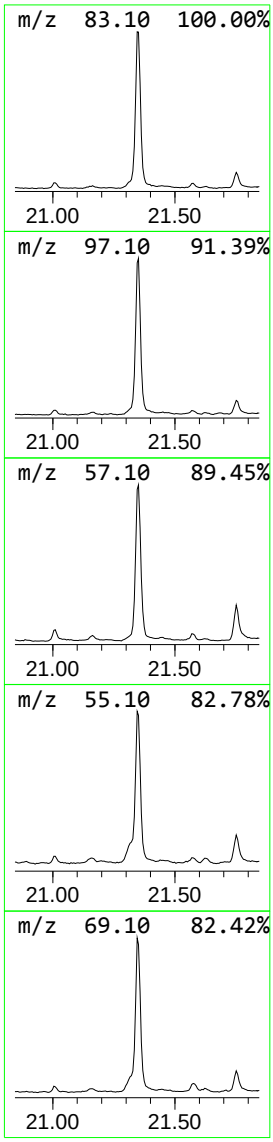
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 n-Nonadecanol-1 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.345	24.24 ng/ul	9212390	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2	Behenic alcohol	326	C22H46O	000661-19-8	94
3	1-Nonadecene	266	C19H38	018435-45-5	91
4	1-Heneicosanol	312	C21H44O	015594-90-8	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91





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Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

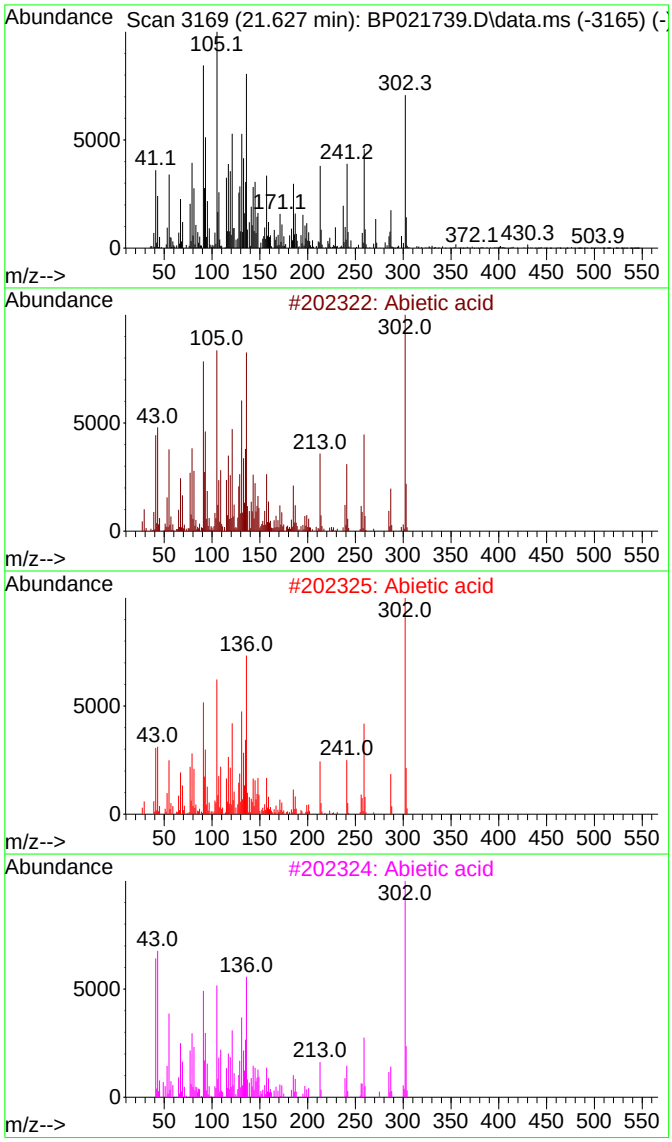
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

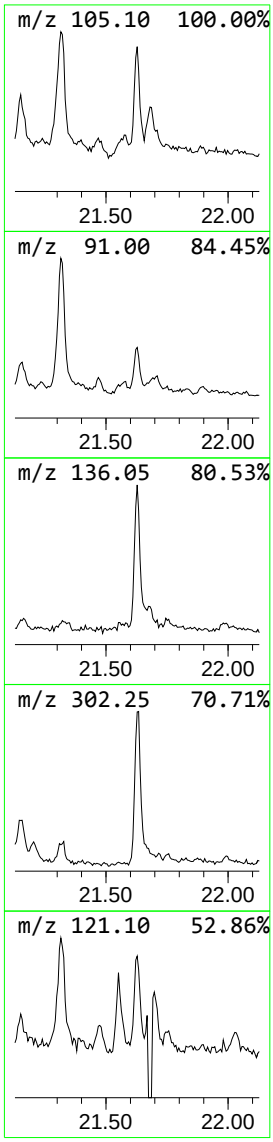
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Abietic acid Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.627	2.55 ng/ul	968016	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Abietic acid	302	C20H30O2	000514-10-3	98
2	Abietic acid	302	C20H30O2	000514-10-3	93
3	Abietic acid	302	C20H30O2	000514-10-3	89
4	Abietic acid	302	C20H30O2	000514-10-3	70
5	Isopimaric acid	302	C20H30O2	005835-26-7	38





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Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

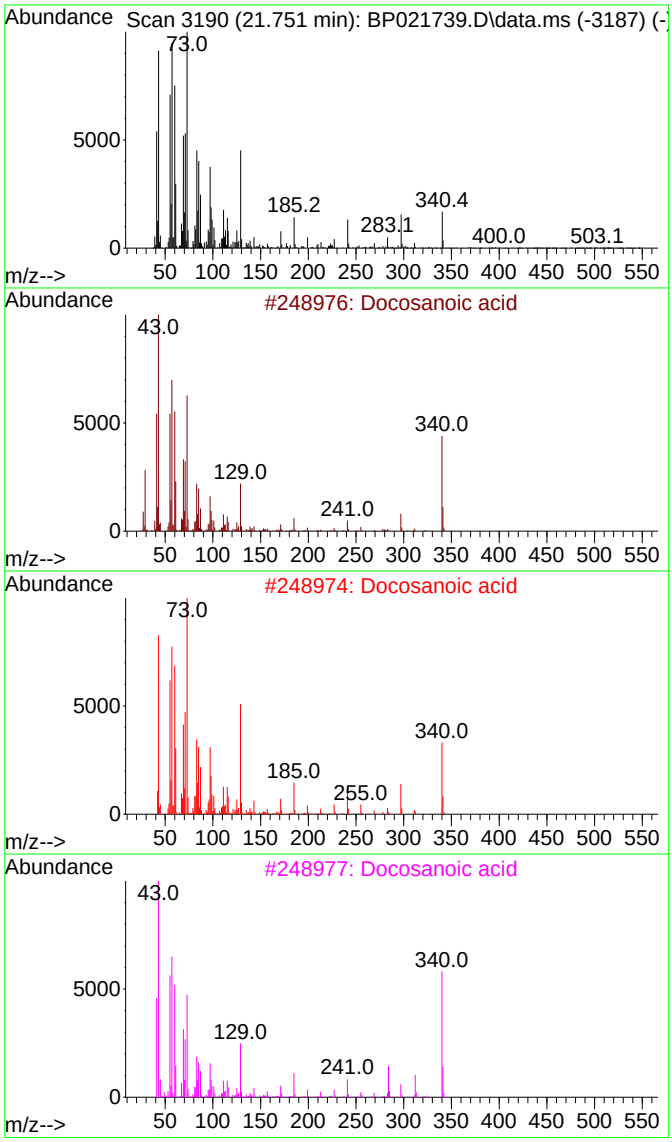
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

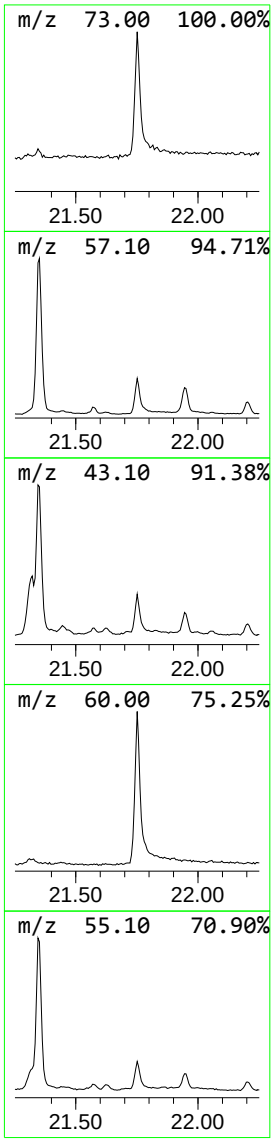
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.751	5.47 ng/ul	2080620	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosanoic acid	340	C22H44O2	000112-85-6	99
2	Docosanoic acid	340	C22H44O2	000112-85-6	96
3	Docosanoic acid	340	C22H44O2	000112-85-6	89
4	Octadecanoic acid	284	C18H36O2	000057-11-4	81
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	70





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Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

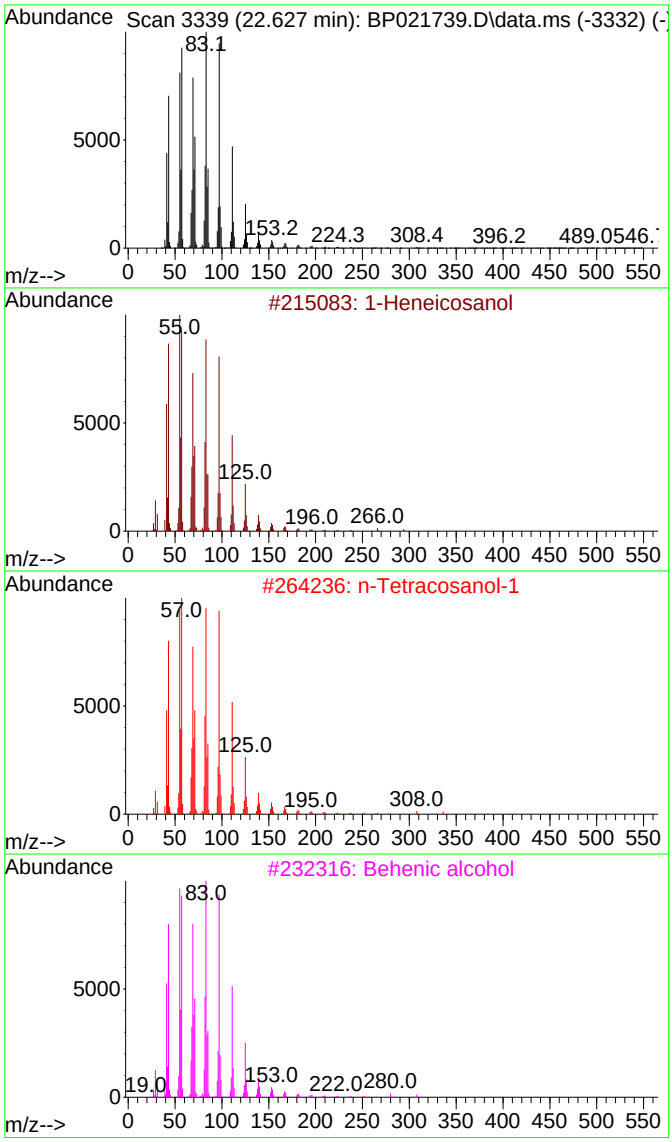
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

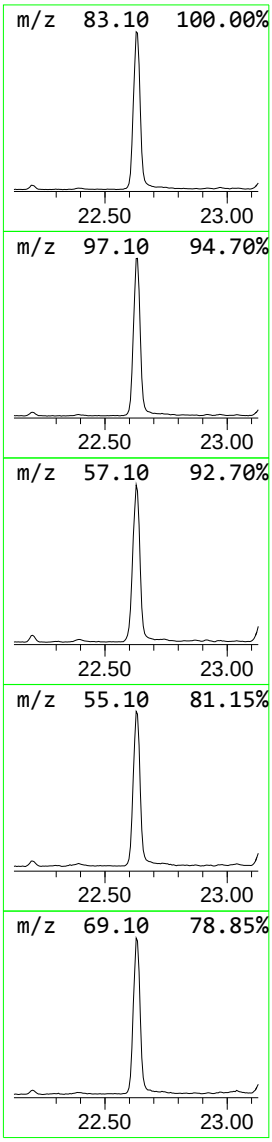
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.627	49.20 ng/ul	18703400	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3	Behenic alcohol	326	C22H46O	000661-19-8	94
4	1-Heptacosanol	396	C27H56O	002004-39-9	93
5	1-Tetracosene	336	C24H48	010192-32-2	91





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Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

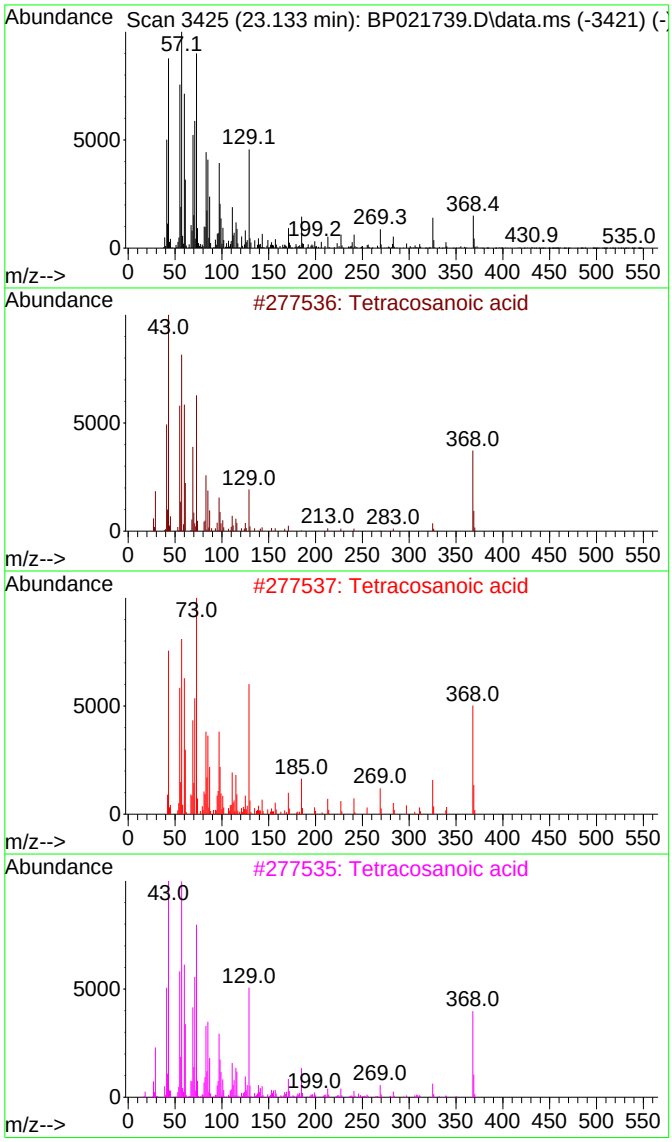
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

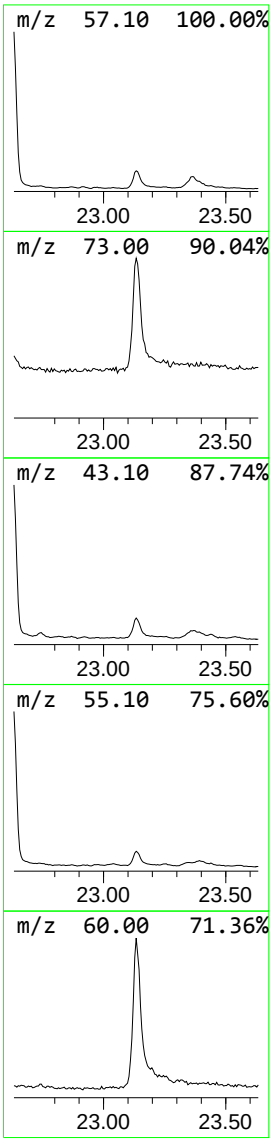
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Tetracosanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.133	6.75 ng/ul	2565520	Chrysene-d12	21.680

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosanoic acid	368	C24H48O2	000557-59-5	99
2	Tetracosanoic acid	368	C24H48O2	000557-59-5	96
3	Tetracosanoic acid	368	C24H48O2	000557-59-5	90
4	Hexacosanoic acid	396	C26H52O2	000506-46-7	83
5	Hexacosyl acetate	424	C28H56O2	000822-32-2	60





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Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

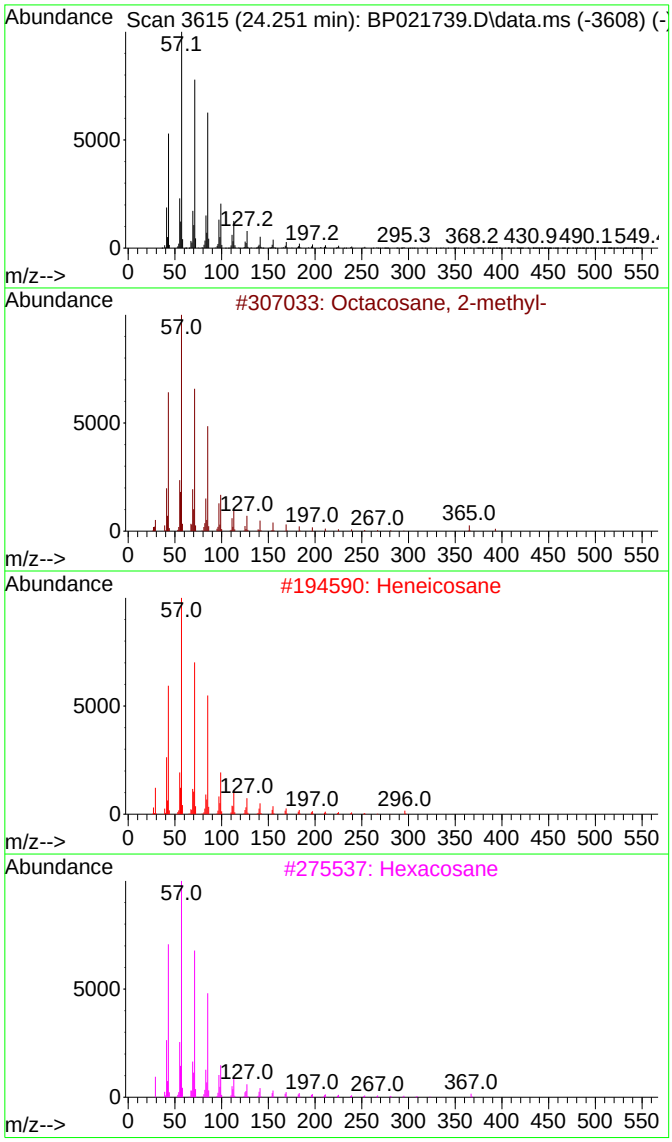
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

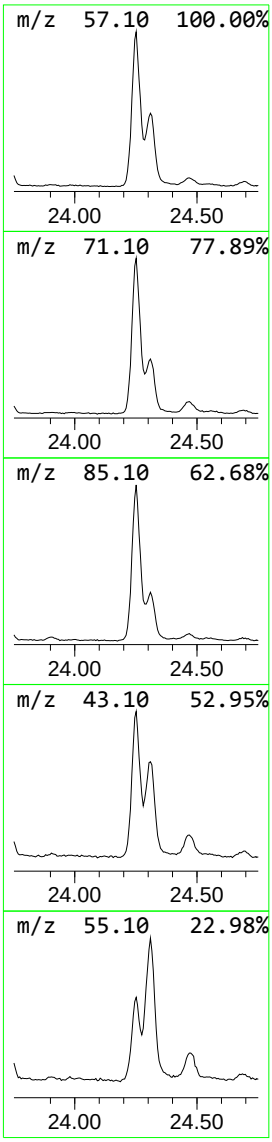
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.251	11.88 ng/ul	4516930	Perylene-d12	25.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

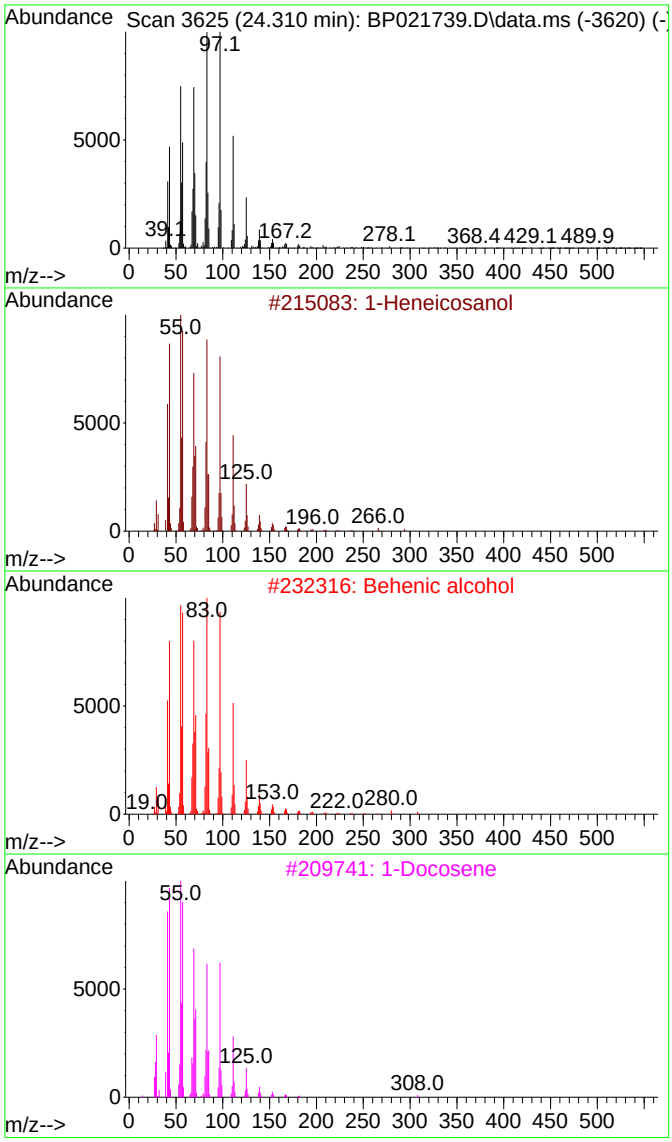
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

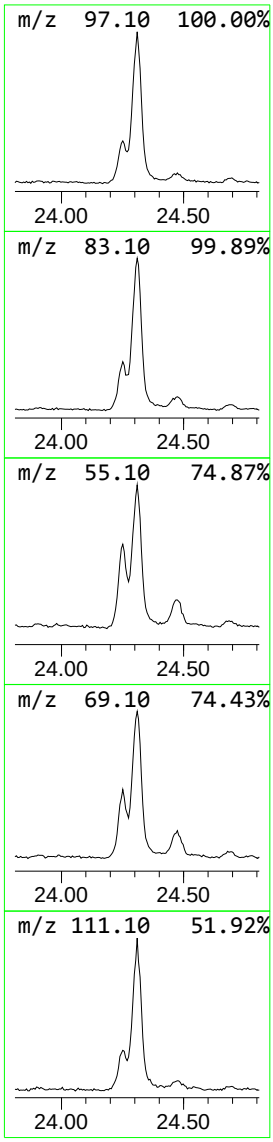
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Behenic alcohol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.310	12.66 ng/ul	4810900	Perylene-d12	25.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	86
2	Behenic alcohol	326	C22H46O	000661-19-8	83
3	1-Docosene	308	C22H44	001599-67-3	64
4	1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	53
5	Oleyl alcohol, trifluoroacetate	364	C20H35F3O2	1000352-68-4	52





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Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

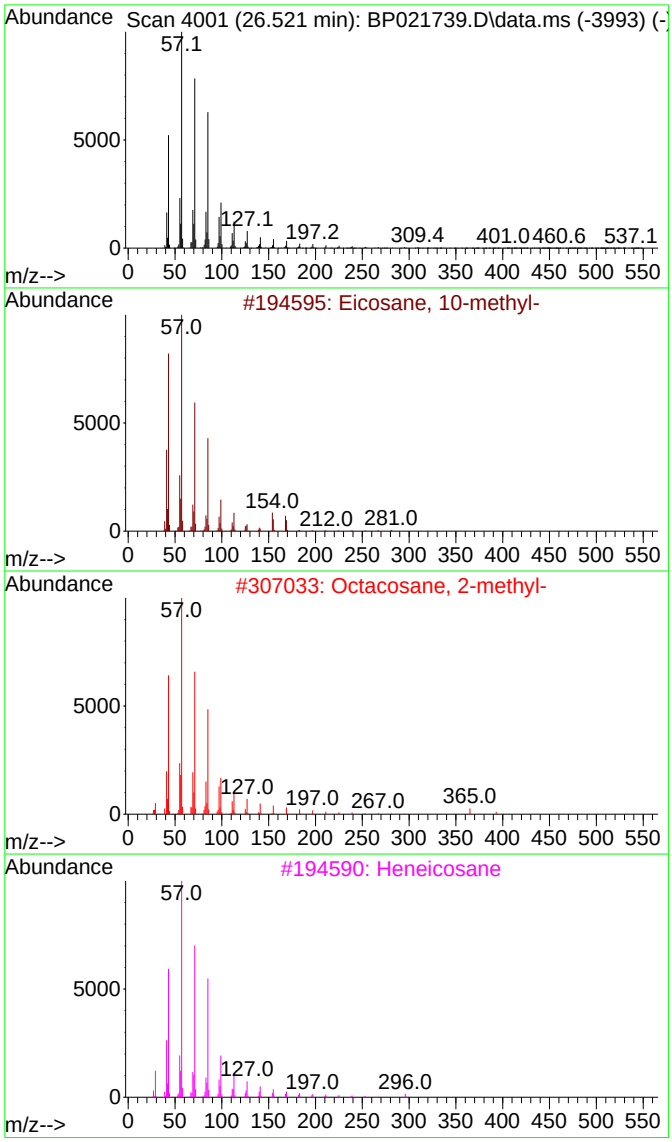
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

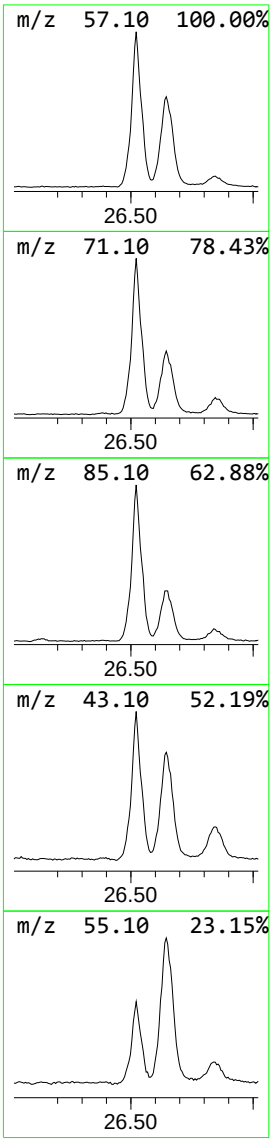
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.521	13.88 ng/ul	5274330	Perylene-d12	25.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
5	Hexacosane	366	C26H54	000630-01-3	90





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Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

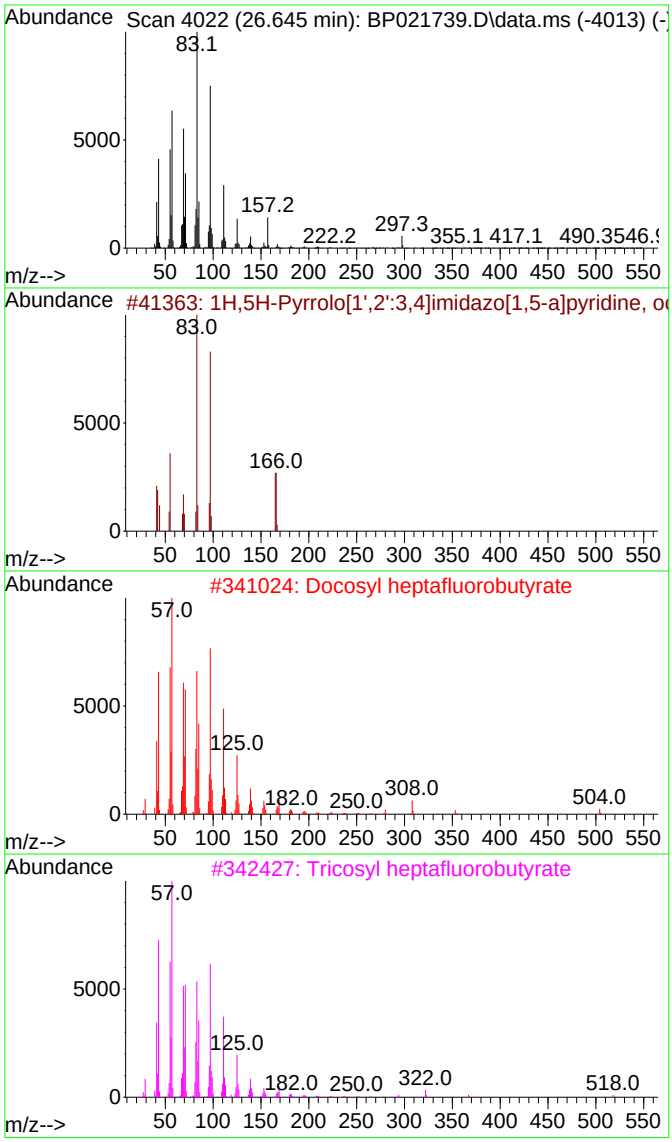
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

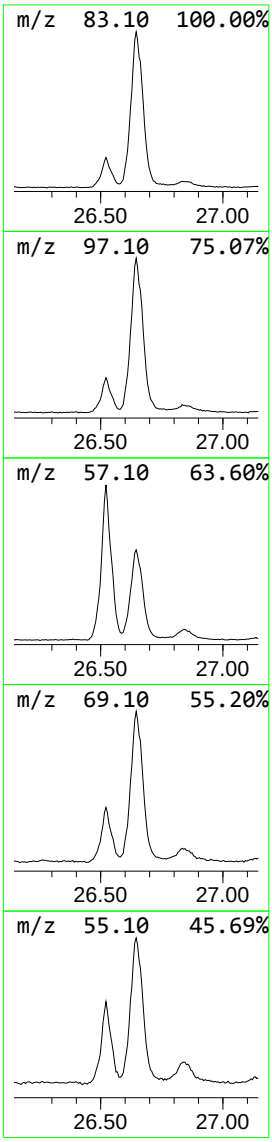
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.645	23.09 ng/ul	8776080	Perylene-d12	25.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H,5H-Pyrrolo[1',2':3,4]imidazo[...	166	C10H18N2	054966-11-9	64
2	Docosyl heptafluorobutyrate	522	C26H45F7O2	1000351-83-1	41
3	Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	41
4	Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	38
5	Triacontan-15-ol	438	C30H62O	031849-26-0	38





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Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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

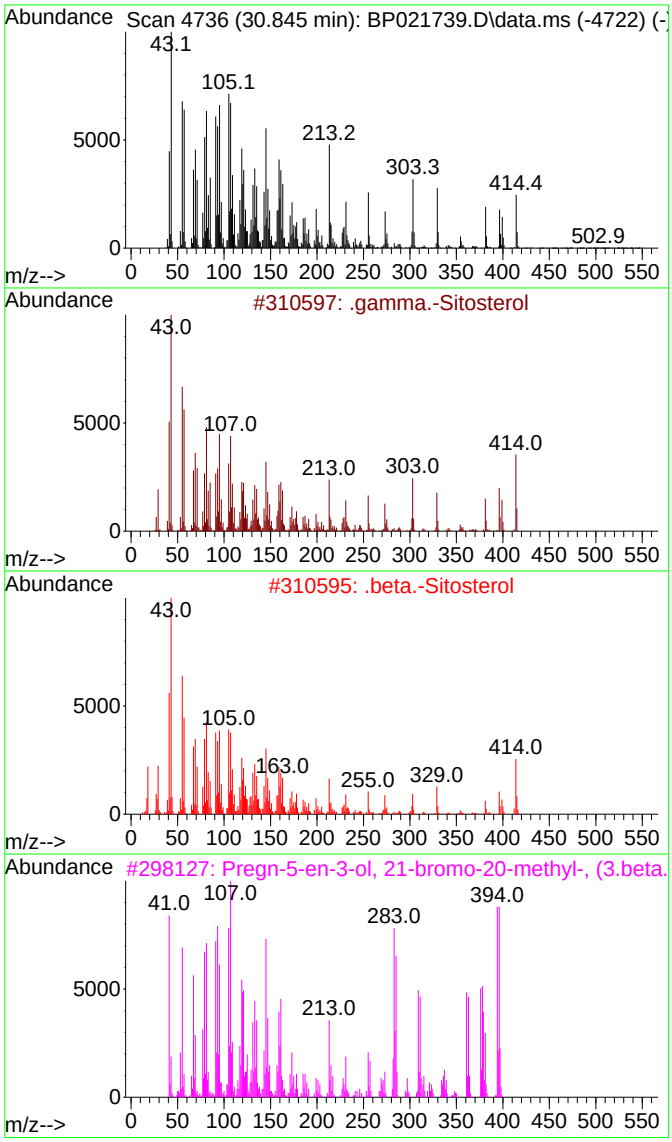
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

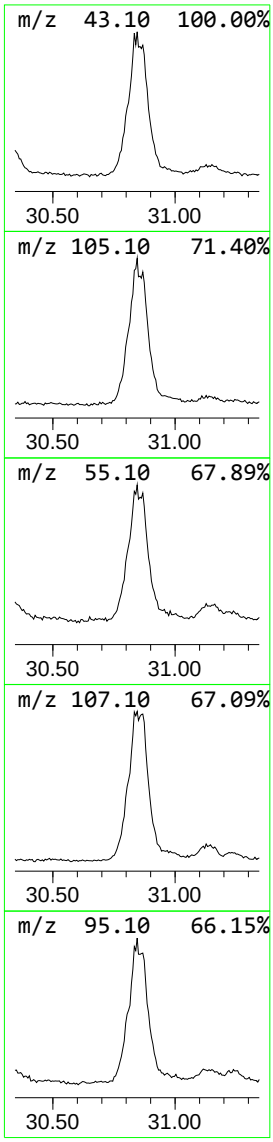
****

Peak Number 25 .gamma.-Sitosterol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.845	21.02 ng/ul	7989960	Perylene-d12	25.098

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	95
3	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	94
4	.beta.-Sitosterol	414	C29H50O	000083-46-5	89
5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021739.D
Acq On : 30 Aug 2024 03:36
Operator : RC/JU
Sample : P3721-13
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ8

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	147.1	ng/ul	23698200	1	7.793	3221530	20.0
(1R)-2,6,6-Trim...	6.499	13.7	ng/ul	2212210	1	7.793	3221530	20.0
Benzeneacetic acid	11.116	2.6	ng/ul	606731	2	10.575	4699160	20.0
n-Hexadecanoic ...	18.163	11.2	ng/ul	4260800	4	17.228	7624490	20.0
m-Camphorene	18.516	2.1	ng/ul	794858	4	17.228	7624490	20.0
unknown-01	19.110	3.8	ng/ul	1428640	4	17.228	7624490	20.0
cis-Vaccenic acid	19.375	15.6	ng/ul	5953910	4	17.228	7624490	20.0
Octadecanoic acid	19.498	4.9	ng/ul	1875530	5	21.680	7602460	20.0
(Z)-5-(Pentadec...	19.622	3.4	ng/ul	1291250	5	21.680	7602460	20.0
(3Z)-1,7,7-trim...	20.016	3.8	ng/ul	1443750	5	21.680	7602460	20.0
Acridin-9-amine...	20.510	4.9	ng/ul	1871740	5	21.680	7602460	20.0
unknown-02	20.592	2.3	ng/ul	863077	5	21.680	7602460	20.0
Methyl dehydroa...	20.733	4.4	ng/ul	1660970	5	21.680	7602460	20.0
Dehydroabietic ...	21.322	39.7	ng/ul	15087700	5	21.680	7602460	20.0
n-Nonadecanol-1	21.345	24.2	ng/ul	9212390	5	21.680	7602460	20.0
Abietic acid	21.627	2.5	ng/ul	968016	5	21.680	7602460	20.0
Docosanoic acid	21.751	5.5	ng/ul	2080620	5	21.680	7602460	20.0
1-Heneicosanol	22.627	49.2	ng/ul	18703400	5	21.680	7602460	20.0
Tetracosanoic acid	23.133	6.8	ng/ul	2565520	5	21.680	7602460	20.0
(DEL) Alkane: S...	24.251	11.9	ng/ul	4516930	6	25.098	7602590	20.0
Behenic alcohol	24.310	12.7	ng/ul	4810900	6	25.098	7602590	20.0
(DEL) Alkane: S...	26.521	13.9	ng/ul	5274330	6	25.098	7602590	20.0
1H,5H-Pyrrolo[1...	26.645	23.1	ng/ul	8776080	6	25.098	7602590	20.0
.gamma.-Sitosterol	30.845	21.0	ng/ul	7989960	6	25.098	7602590	20.0