LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034994.D
Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

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_M\Methods\SFAM-EPA-BM051022.M
Title : SVOA CALIBRATION

Signal : TIC: BM034994.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	44	48	58	rVB	72808	103552	1.86%	0.135%
2	3.781	114	118	132	rBV	503159	767610	13.78%	1.000%
3	3.887	132	136	144	rVV	102822	141277	2.54%	0.184%
4	4.016	156	158	164	rVB	15139	22757	0.41%	0.030%
5	4.110	171	174	179	rVB2	21615	30492	0.55%	0.040%
6	4.669	264	269	273	rBV4	16942	24983	0.45%	0.033%
7	5.028	325	330	337	rVB	58899	88174	1.58%	0.115%
8	5.122	342	346	352	rVB2	18673	29297	0.53%	0.038%
9	6.910	643	650	657	rBV6	13329	23895	0.43%	0.031%
10	7.087	670	680	697	rVB	1161523	1960482	35.18%	2.554%
11	7.251	701	708	718	rBV	941932	1462208	26.24%	1.905%
12	7.451	734	742	754	rBV	1239092	1985146	35.63%	2.586%
13	7.557	756	760	774	rVV8	19503	60710	1.09%	0.079%
14	7.922	815	822	830	rVV	791151	1243476	22.32%	1.620%
15	8.157	855	862	871	rBV	227496	370980	6.66%	0.483%
16	8.628	935	942	950	rBV	1068442	1693977	30.40%	2.207%
17	8.745	958	962	968	rVB	13180	25984	0.47%	0.034%
18	9.081	1010	1019	1026	rVV	1145802	1901953	34.13%	2.477%
19	9.257	1043	1049	1054	rBV4	16787	29387	0.53%	0.038%
20	9.804	1135	1142	1149	rVV	968987	1591278	28.56%	2.073%
21	10.339	1226	1233	1243	rBV	1282890	2180699	39.14%	2.840%
22	10.428	1243	1248	1254	rVV4	16587	34480	0.62%	0.045%
23	10.504	1254	1261	1272	rVV	122613	199226	3.58%	0.260%
24	10.722	1291	1298	1305	rBV	1001905	1700783	30.52%	2.215%
25	10.769	1305	1306	1313	rVB	23940	33009	0.59%	0.043%
26	10.857	1313	1321	1327	rBV	942731	1631357	29.28%	2.125%
27	11.692	1450	1463	1470	rBV	30142	65812	1.18%	0.086%
28	11.763	1470	1475	1481	rBV2	17141	31699	0.57%	0.041%
29	12.098	1527	1532	1537	rVV	53764	85871	1.54%	0.112%
30	12.163	1537	1543	1549	rVV	33130	54279	0.97%	0.071%
31	12.310	1560	1568	1572	rVV	28180	57116	1.03%	0.074%
32	12.380	1576	1580	1590	rVB2	23743	41639	0.75%	0.054%
33	12.551	1600	1609	1614	rBV4	8019	25225	0.45%	0.033%
34	12.922	1669	1672	1677	rVV7	14072	24895	0.45%	0.032%
35	13.074	1690	1698	1700	rBV8	11041	26046	0.47%	0.034%
36	13.110	1701	1704	1710	rVV3	26179	35238	0.63%	0.046%

LSC Area Percent Report

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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : EW497

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_M\Methods\SFAM-EPA-BM051022.M
Title : SVOA CALIBRATION

37	13.174	1710	1715	1722	rVB4	17576	27369	0.49%	0.036%
38	13.398	1748	1753	1758	rVB2	50747	74134	1.33%	0.097%
39	13.469	1760	1765	1769	rBV6	16041	30403	0.55%	0.040%
40	13.874	1830	1834	1838	rVV6	16532	26945	0.48%	0.035%
41	13.963	1843	1849	1855	rVB	2105239	2817111	50.56%	3.669%
42	14.051	1861	1864	1868	rVB2	27038	33371	0.60%	0.043%
43	14.098	1868	1872	1878	rVB8	17981	29388	0.53%	0.038%
44	14.245	1887	1897	1907	rBV	2365728	3583145	64.30%	4.667%
45	14.392	1919	1922	1930	rVB8	12924	26368	0.47%	0.034%
46	14.463	1930	1934	1940	rVB	73056	95212	1.71%	0.124%
47	14.551	1940	1949	1957	rBV	1264512	1975578	35.45%	2.573%
48	14.616	1957	1960	1969	rVB2	29118	50286	0.90%	0.066%
49	14.739	1975	1981	1996	rBV	741024	1164984	20.91%	1.517%
50	14.892	2003	2007	2012	rBV7	38176	71262	1.28%	0.093%
51	14.963	2014	2019	2025	rVB3	47115	94572	1.70%	0.123%
52	15.086	2036	2040	2045	rVB3	27795	44271	0.79%	0.058%
53	15.363	2082	2087	2092	rBV3	42600	79975	1.44%	0.104%
54	15.416	2092	2096	2100	rVB	104645	125090	2.24%	0.163%
55	15.457	2100	2103	2110	rBV4	14821	25458	0.46%	0.033%
56	15.545	2112	2118	2124	rVV	2946272	4097121	73.53%	5.337%
57	15.598	2124	2127	2130	rVV4	39235	47852	0.86%	0.062%
58	15.657	2132	2137	2144	rVB	724456	963800	17.30%	1.255%
59	15.786	2156	2159	2161	rBV	25310	28894	0.52%	0.038%
60	15.821	2163	2165	2168	rVB	40335	38875	0.70%	0.051%
61	15.915	2177	2181	2187	rVB3	48611	78314	1.41%	0.102%
62	15.992	2187	2194	2198	rBV2	51717	117981	2.12%	0.154%
63	16.033	2198	2201	2204	rVB4	27164	35245	0.63%	0.046%
64	16.245	2231	2237	2240	rBV	476706	645947	11.59%	0.841%
65	16.298	2240	2246	2250	rVB2	267050	484709	8.70%	0.631%
66	16.386	2256	2261	2265	rBV	413360	557163	10.00%	0.726%
67	16.445	2265	2271	2277	rVV3	523698	999901	17.94%	1.302%
68	16.504	2277	2281	2285	rVV2	544266	761886	13.67%	0.992%
69	16.551	2285	2289	2293	rVV	291039	401486	7.21%	0.523%
70	16.604	2293	2298	2299	rVV	134303	200374	3.60%	0.261%
71	16.621	2299	2301	2304	rVV	184297	264217	4.74%	0.344%
72	16.651	2304	2306	2309	rVV	193420	211950	3.80%	0.276%
73	16.704	2309	2315	2322	rVV4	369252	761063	13.66%	0.991%
74	16.768	2322	2326	2330	rVV	489181	686371	12.32%	0.894%
75	16.810	2330	2333	2335	rVV	161870	230754	4.14%	0.301%
76	16.845	2335	2339	2347	rVB2	301388	450906	8.09%	0.587%

LSC Area Percent Report

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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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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
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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Title : SVOA CALIBRATION

77	17.068	2373	2377	2380	rBV	115172	139756	2.51%	0.182%
78	17.292	2410	2415	2420	rBV	1297636	1909164	34.26%	2.487%
79	17.333	2420	2422	2426	rVV	303769	378300	6.79%	0.493%
80	17.392	2426	2432	2444	rVB	2762792	3925959	70.46%	5.114%
81	17.804	2499	2502	2507	rVB	89637	111135	1.99%	0.145%
82	18.192	2565	2568	2576	rVB2	261371	385811	6.92%	0.503%
83	18.498	2617	2620	2629	rVB2	148672	199827	3.59%	0.260%
84	19.127	2723	2727	2734	rBV	356189	428884	7.70%	0.559%
85	19.345	2760	2764	2771	rVB	520602	686376	12.32%	0.894%
86	19.468	2782	2785	2788	rBV	171513	191710	3.44%	0.250%
87	19.680	2816	2821	2833	rBV2	3120013	5572182	100.00%	7.258%
88	20.239	2912	2916	2920	rVB	790256	856938	15.38%	1.116%
89	20.733	2996	3000	3003	rBV	847018	904899	16.24%	1.179%
90	21.192	3074	3078	3085	rVB	1270458	1597165	28.66%	2.080%
91	21.468	3120	3125	3129	rBV	1543165	2163237	38.82%	2.818%
92	21.633	3150	3153	3158	rVB	928503	993412	17.83%	1.294%
93	22.097	3228	3232	3243	rVB	914648	1398036	25.09%	1.821%
94	22.592	3312	3316	3323	rBV	736088	1028158	18.45%	1.339%
95	23.144	3405	3410	3414	rBV2	917721	1443581	25.91%	1.880%
96	23.703	3499	3505	3511	rBV2	2257501	4234304	75.99%	5.515%
97	23.768	3512	3516	3523	rVB2	728480	1332150	23.91%	1.735%
98	23.856	3525	3531	3537	rVB2	1029808	1959332	35.16%	2.552%
99	24.491	3634	3639	3647	rVB	802776	1606328	28.83%	2.092%
100	25.327	3777	3781	3791	rVB	517429	1100675	19.75%	1.434%

Sum of corrected areas: 76772042

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034994.D
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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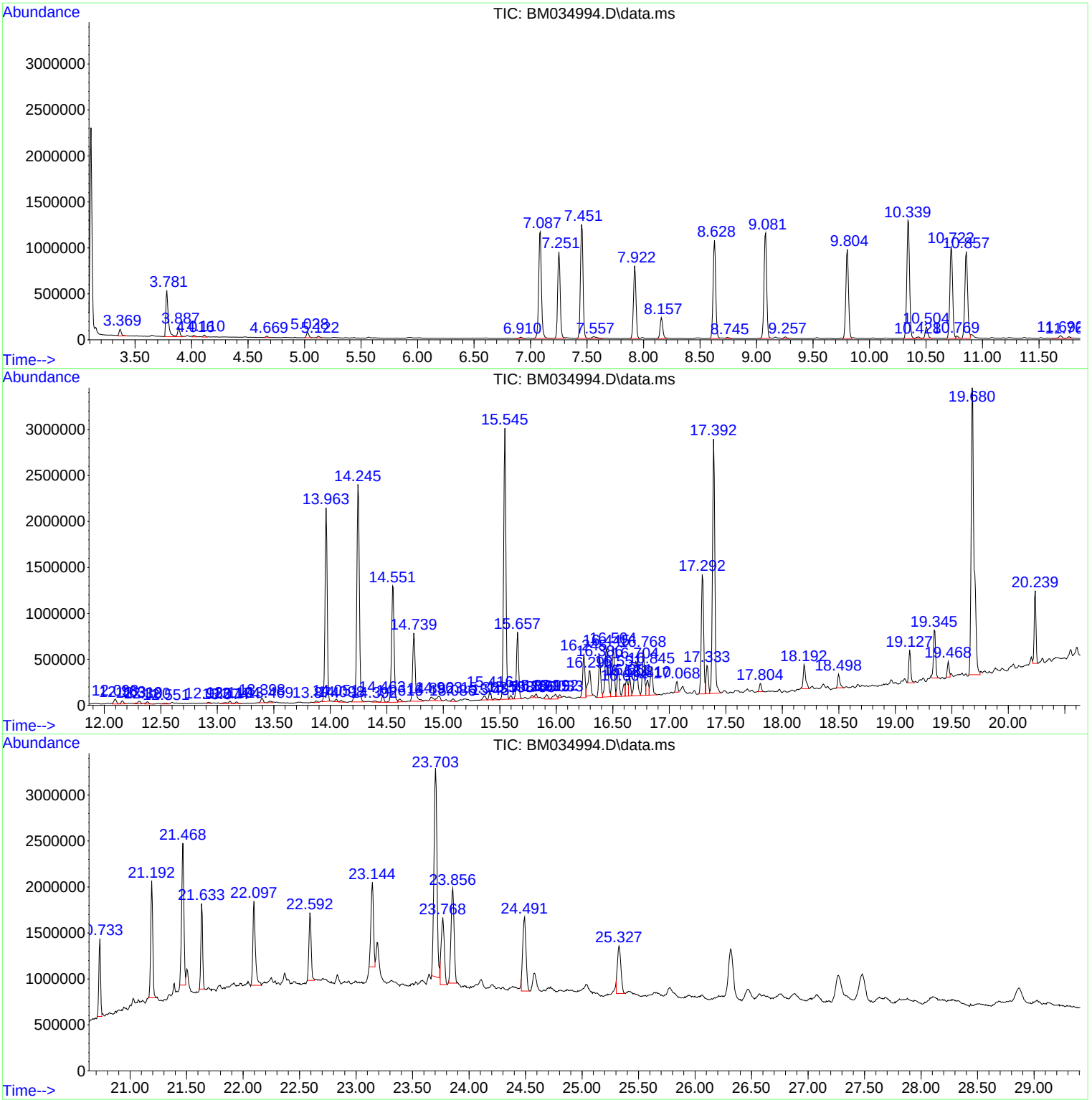
ClientSampleId :
EW497

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034994.D
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

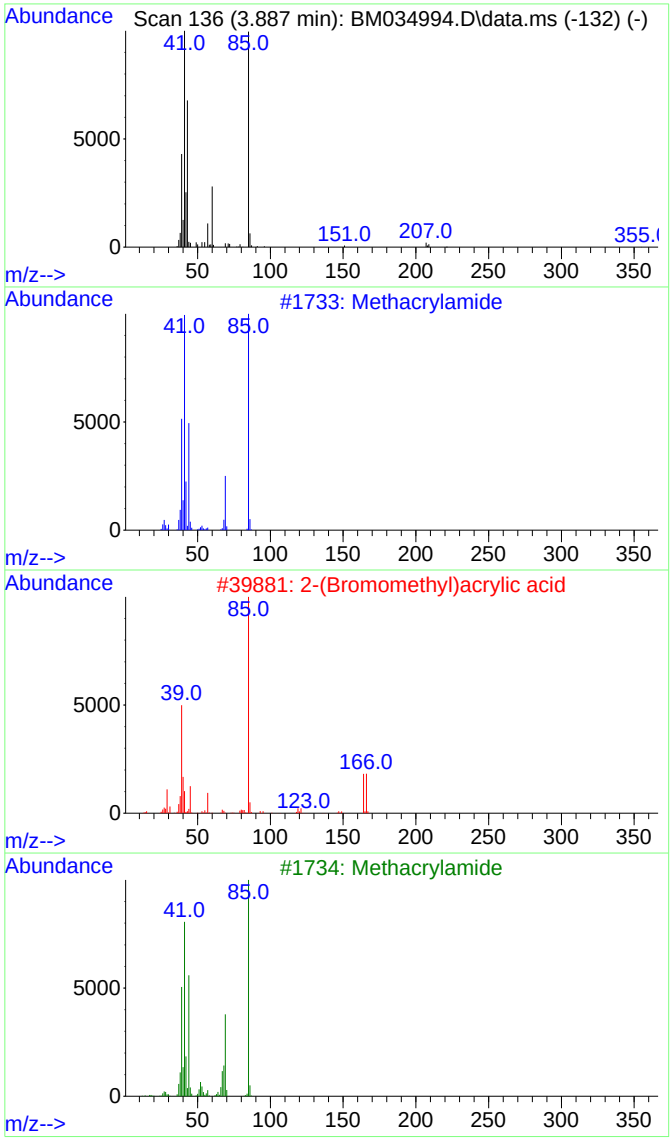
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

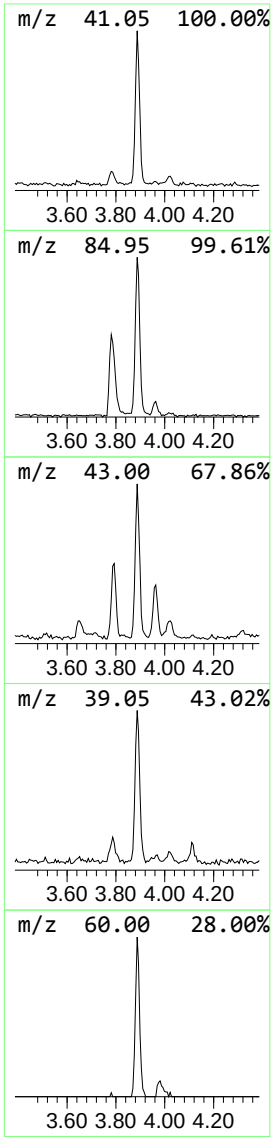
****

Peak Number 1 Methacrylamide Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.887	2.27 ng/ul	141277	1,4-Dichlorobenzene-d4	7.922

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methacrylamide	85	C4H7NO	000079-39-0	50
2	2-(Bromomethyl)acrylic acid	164	C4H5BrO2	072707-66-5	40
3	Methacrylamide	85	C4H7NO	000079-39-0	38
4	Methacrylamide	85	C4H7NO	000079-39-0	28
5	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034994.D
Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

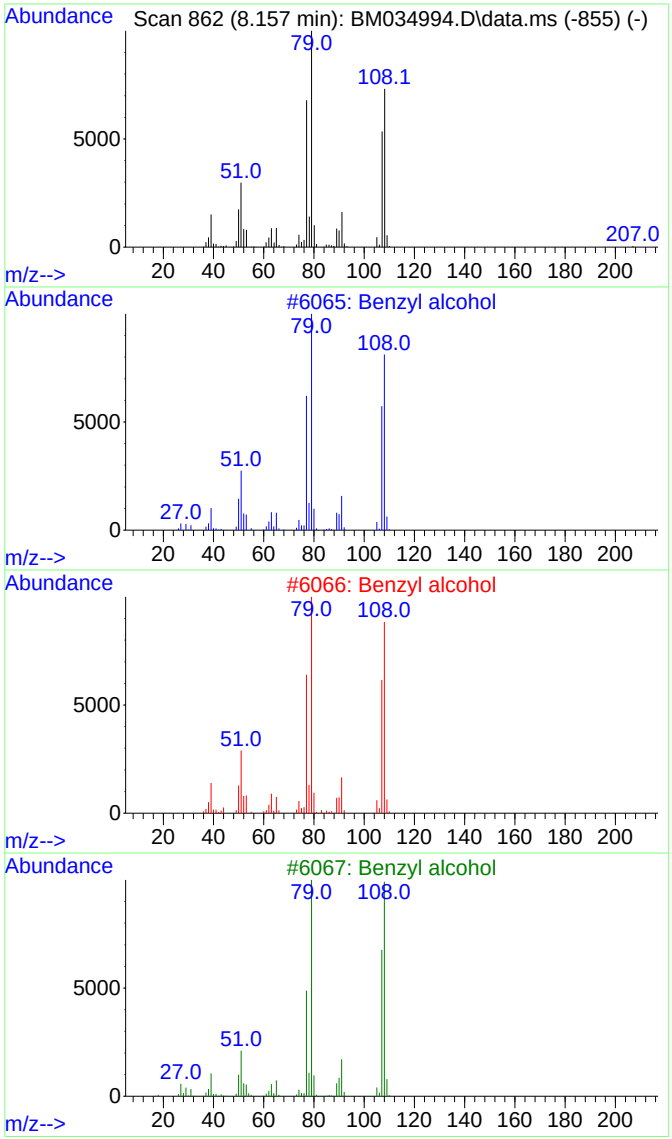
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

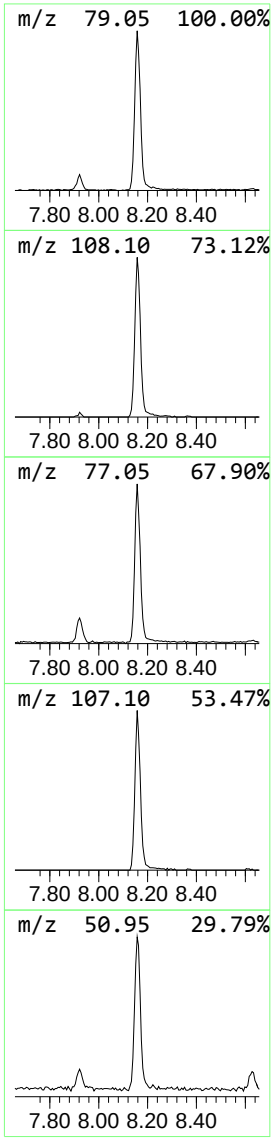
****

Peak Number 2 Benzyl alcohol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.157	5.97 ng/ul	370980	1,4-Dichlorobenzene-d4	7.922

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzyl alcohol	108	C7H8O	000100-51-6	98
2	Benzyl alcohol	108	C7H8O	000100-51-6	97
3	Benzyl alcohol	108	C7H8O	000100-51-6	96
4	Benzyl alcohol	108	C7H8O	000100-51-6	95
5	N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
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ALS Vial : 19 Sample Multiplier: 1

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BNA_M
ClientSampleId :
EW497

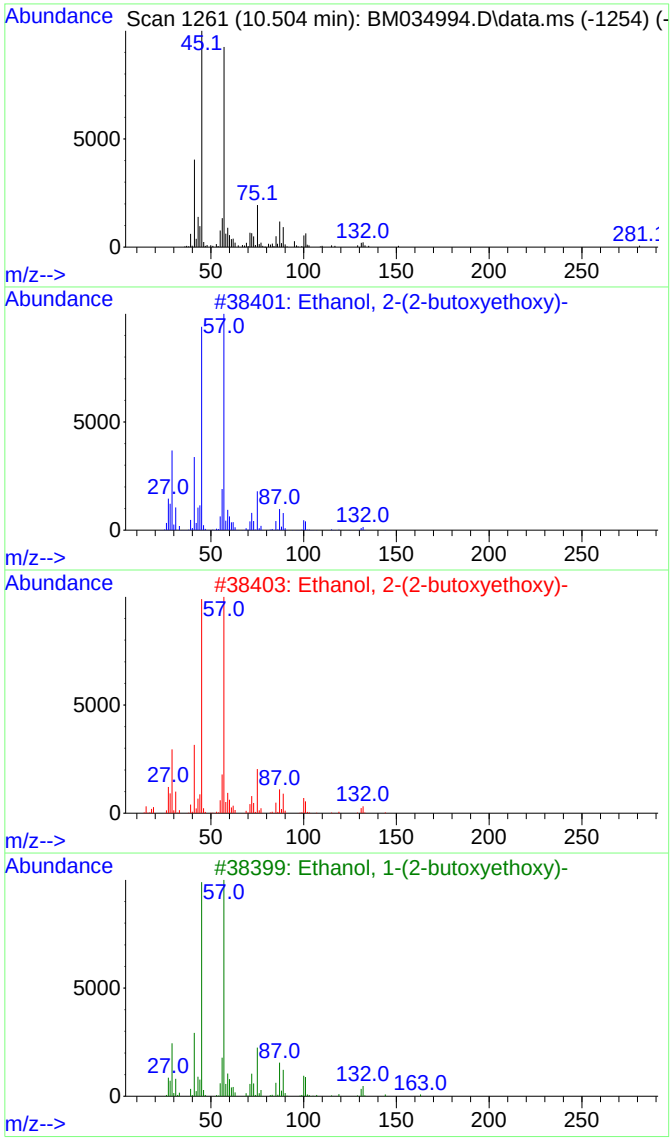
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

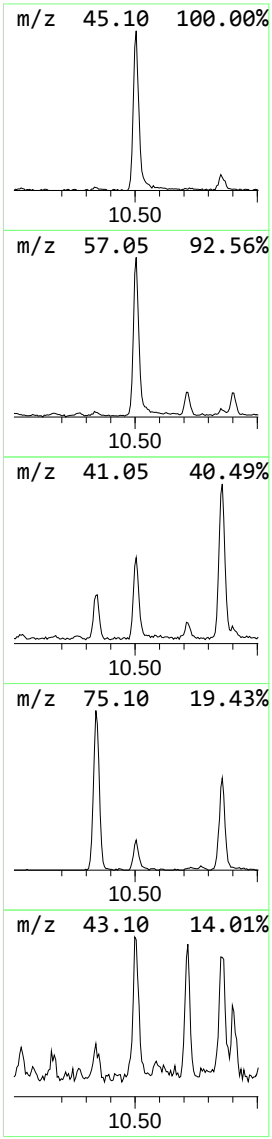
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.504	2.34 ng/ul	199226	Naphthalene-d8	10.722

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
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ALS Vial : 19 Sample Multiplier: 1

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

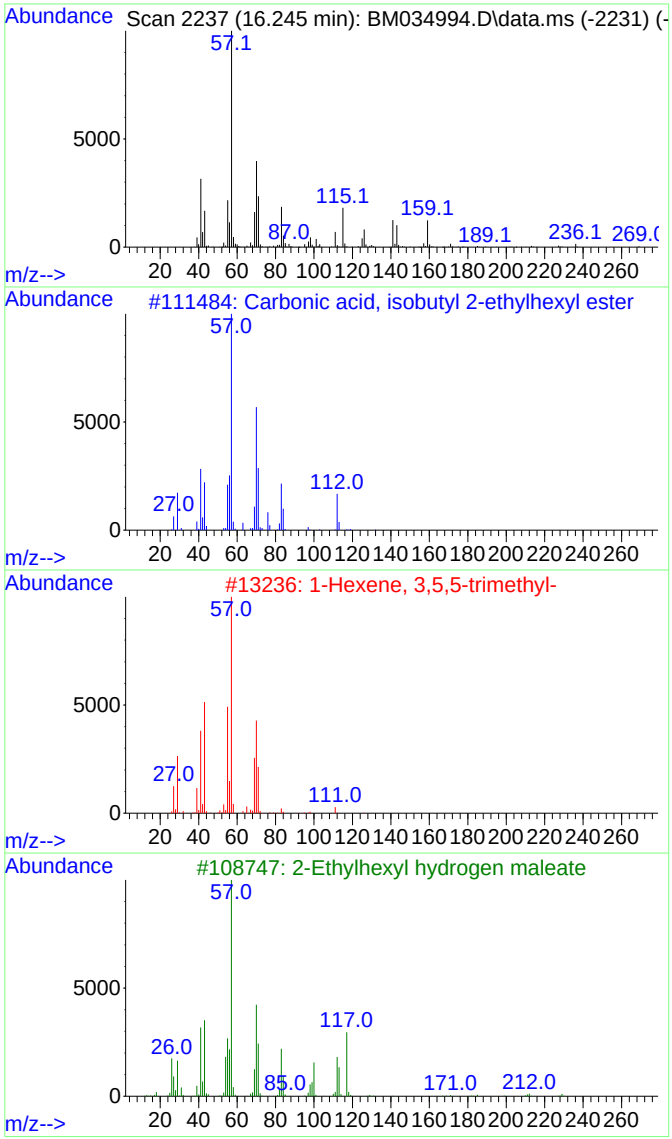
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

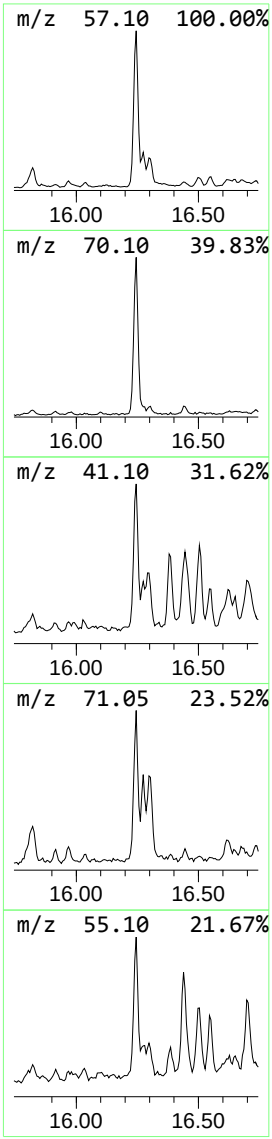
****

Peak Number 4 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.245	6.77 ng/ul	645947	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000383-13-3	47
2	1-Hexene, 3,5,5-trimethyl-	126	C9H18	004316-65-8	38
3	2-Ethylhexyl hydrogen maleate	228	C12H20O4	002370-71-0	38
4	Propanoic acid, 2,2-dimethyl-, 2...	214	C13H26O2	016387-18-1	38
5	Hexanoic acid, 3,5,5-trimethyl-,...	270	C17H34O2	1000406-82-2	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034994.D
Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

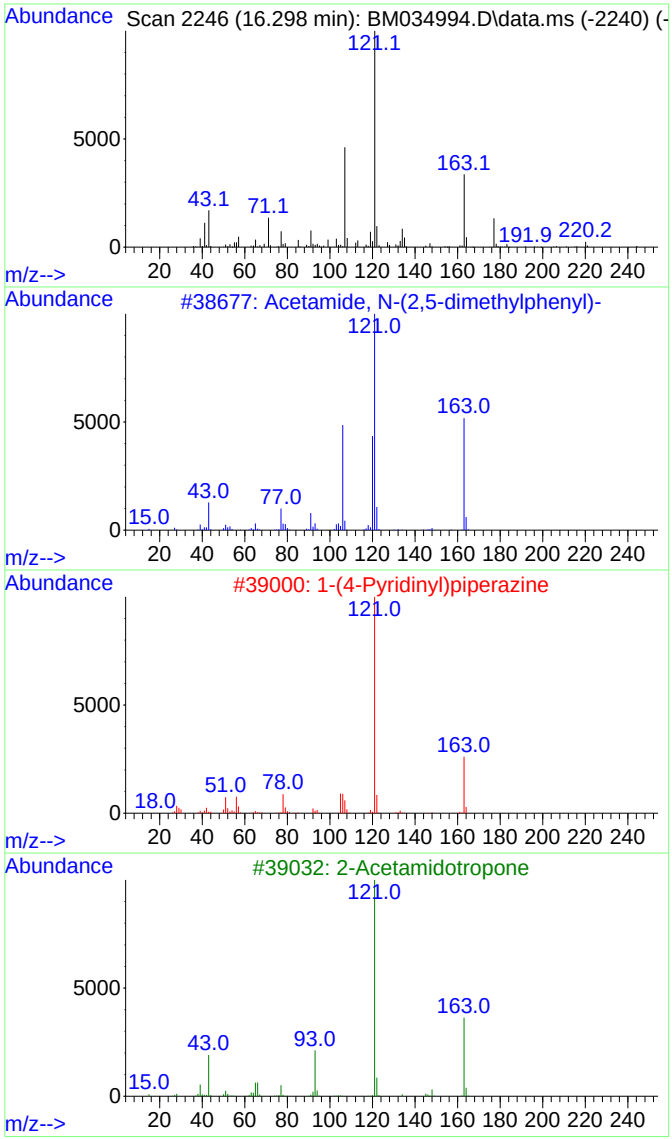
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

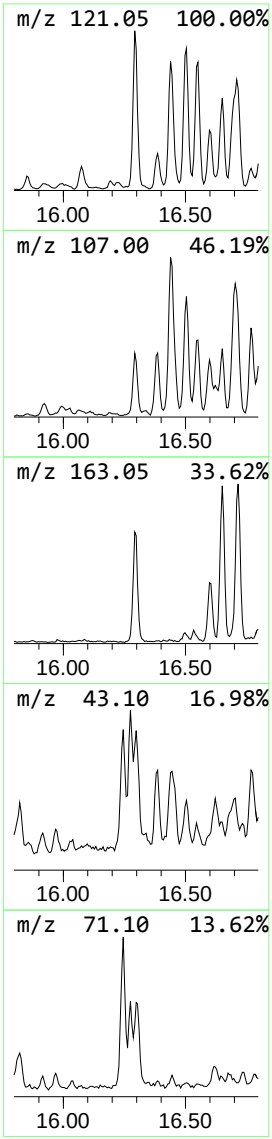
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Acetamide, N-(2,5-dimethylp... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.298	5.08 ng/ul	484709	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2,5-dimethylphenyl)-	163	C10H13NO	002050-44-4	50
2	1-(4-Pyridinyl)piperazine	163	C9H13N3	001008-91-9	50
3	2-Acetamidotropone	163	C9H9NO2	006422-12-4	47
4	N-Allyl-p-hydroxybenzamide	177	C10H11NO2	090921-72-5	46
5	4-sec-Butylphenol, O-isopropylox...	236	C14H20O3	1201410-83-4	43





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Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

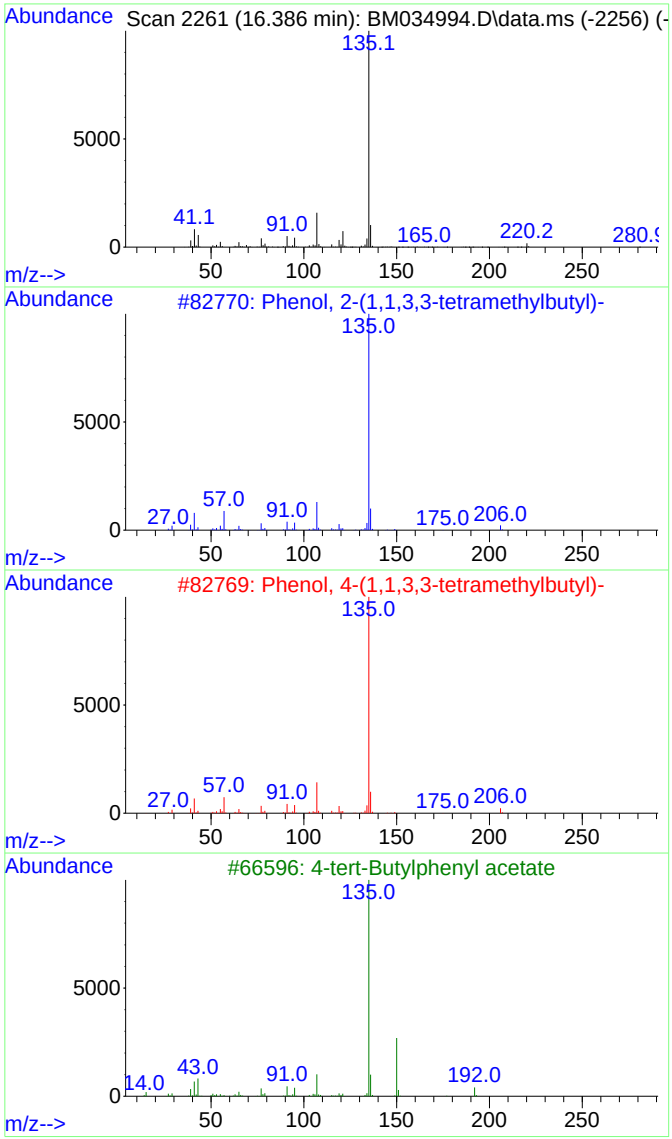
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

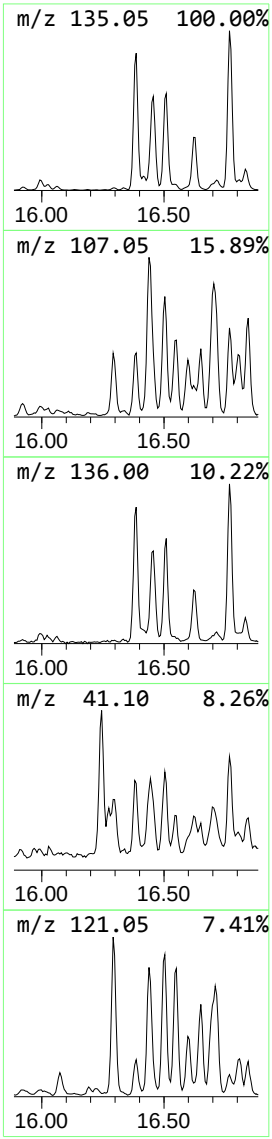
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenol, 2-(1,1,3,3-tetramet... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.386	5.84 ng/ul	557163	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	86
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
3	4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78
4	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
5	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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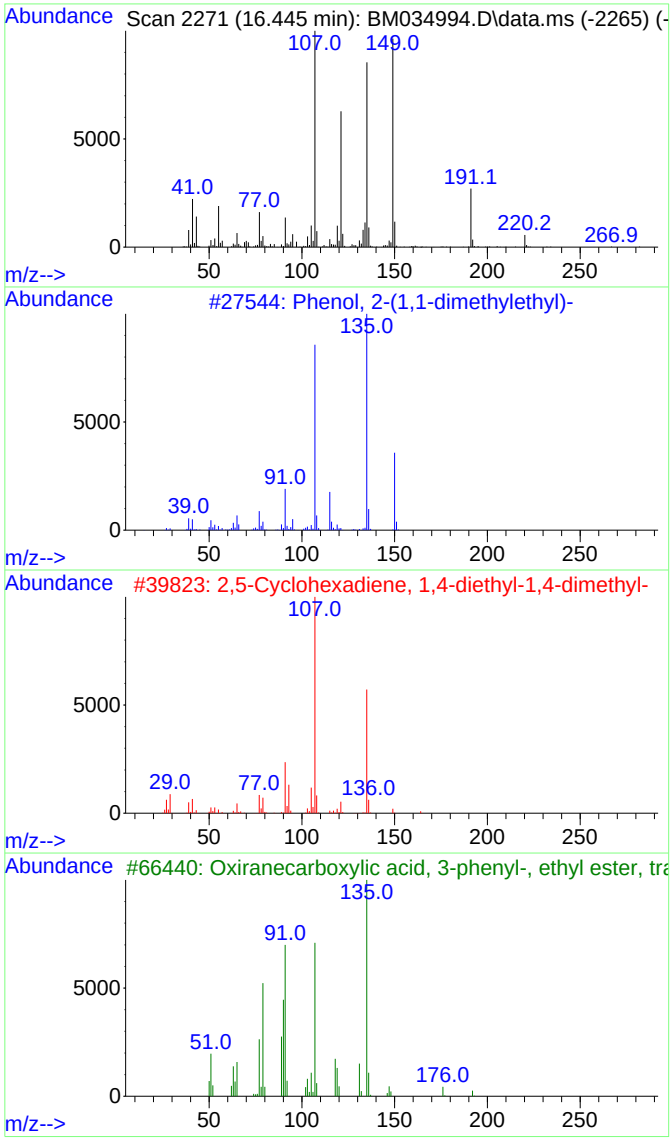
Quant Title : SVOA CALIBRATION

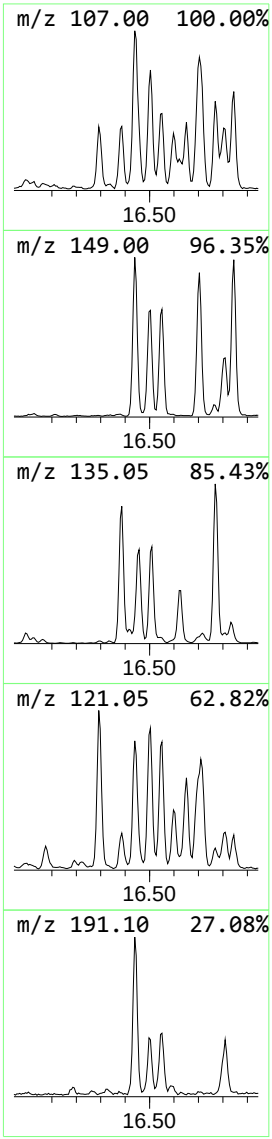
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.445	10.47 ng/ul	999901	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	46
2	2,5-Cyclohexadiene, 1,4-diethyl-...	164	C12H20	1000150-21-6	43
3	Oxiranecarboxylic acid, 3-phenyl...	192	C11H12O3	002272-55-1	38
4	1-Benzoyl-2-piperonyl hydrazine	270	C15H14N2O3	1000256-61-5	35
5	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	35





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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

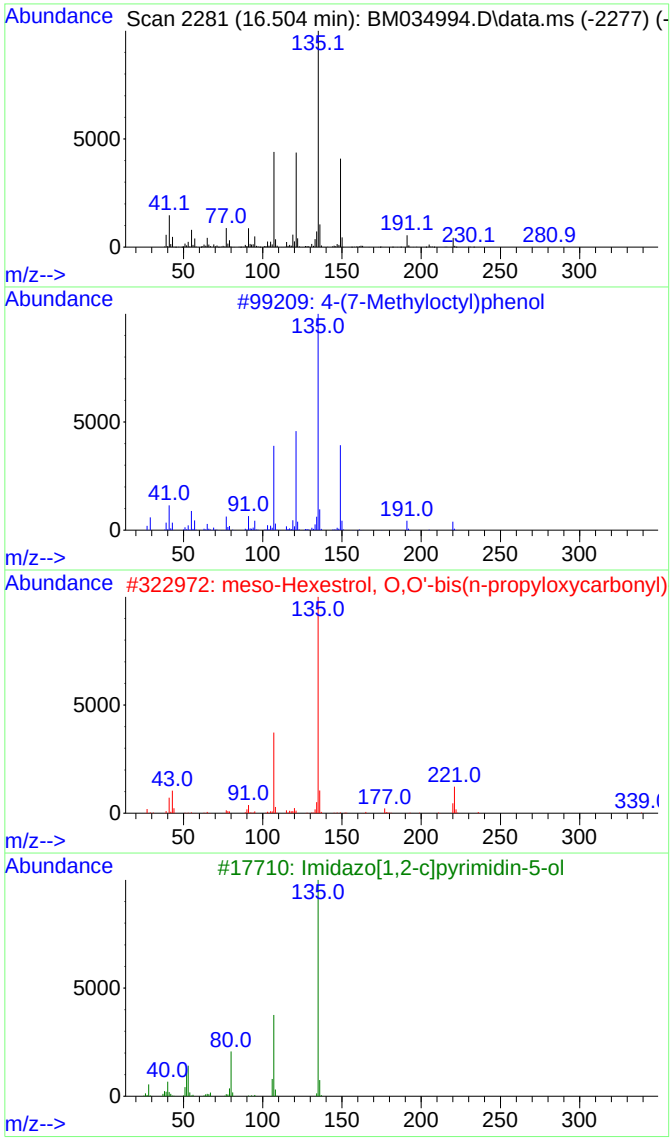
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

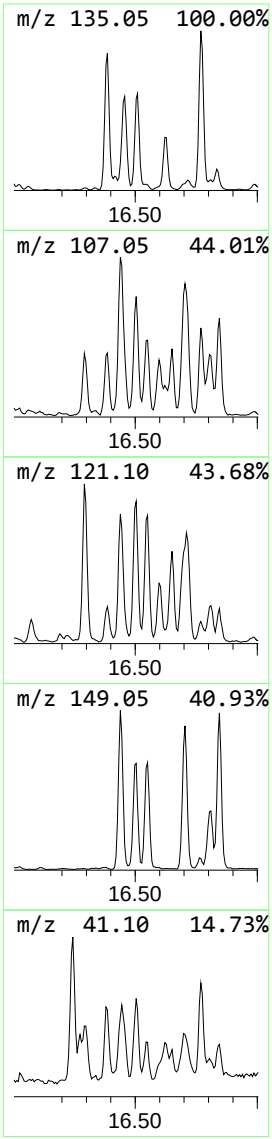
****

Peak Number 8 4-(7-Methyloctyl)phenol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.504	7.98 ng/ul	761886	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	98
2	meso-Hexestrol, O,O'-bis(n-propy...	442	C26H34O6	1000487-65-0	53
3	Imidazo[1,2-c]pyrimidin-5-ol	135	C6H5N3O	055662-66-3	50
4	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	50
5	Hexestrol, O-heptafluorobutyryl-	466	C22H21F7O3	1000365-31-9	50





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Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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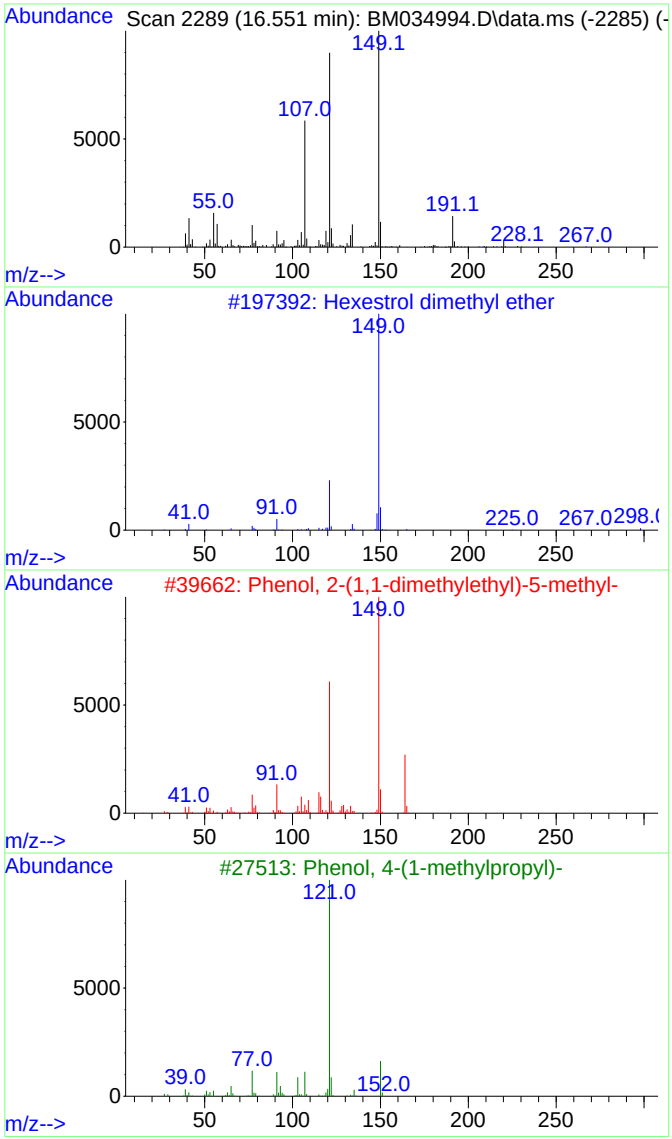
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

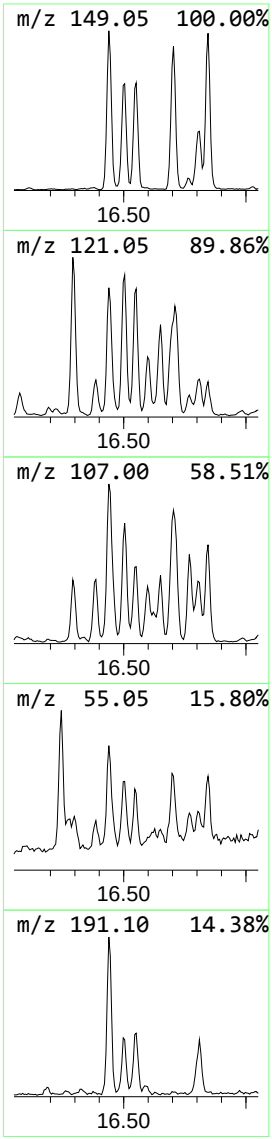
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Hexestrol dimethyl ether Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.551	4.21 ng/ul	401486	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexestrol dimethyl ether	298	C20H26O2	000130-78-9	59
2	Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	59
3	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	30
4	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	30
5	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	27





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Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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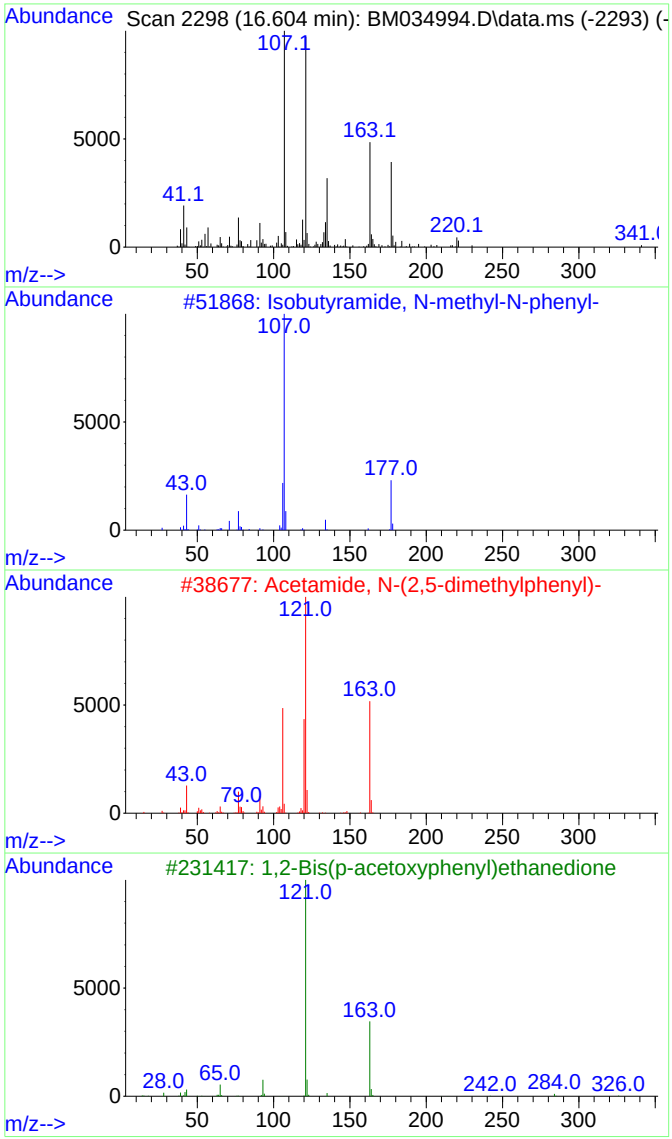
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

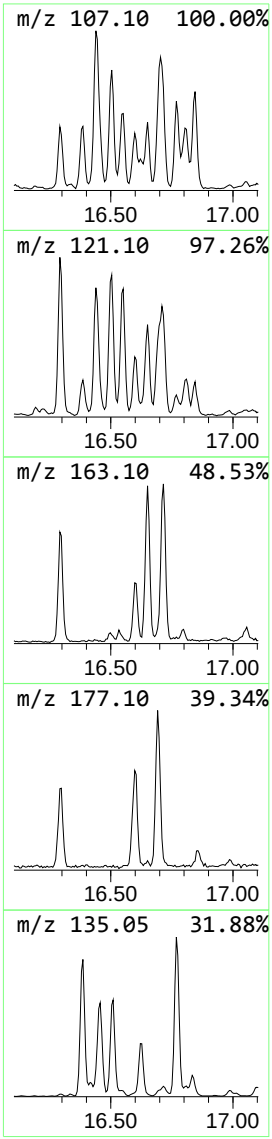
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 unknown-03 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.604	2.10 ng/ul	200374	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isobutyramide, N-methyl-N-phenyl-	177	C11H15NO	1000339-96-1	35
2	Acetamide, N-(2,5-dimethylphenyl)-	163	C10H13NO	002050-44-4	30
3	1,2-Bis(p-acetoxyphenyl)ethanedione	326	C18H14O6	093655-79-9	27
4	N-Cyclopropyl-2-hydroxybenzamide	177	C10H11NO2	440111-82-0	27
5	Butanamide, N-(3-methylphenyl)-	177	C11H15NO	1000306-91-3	27





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Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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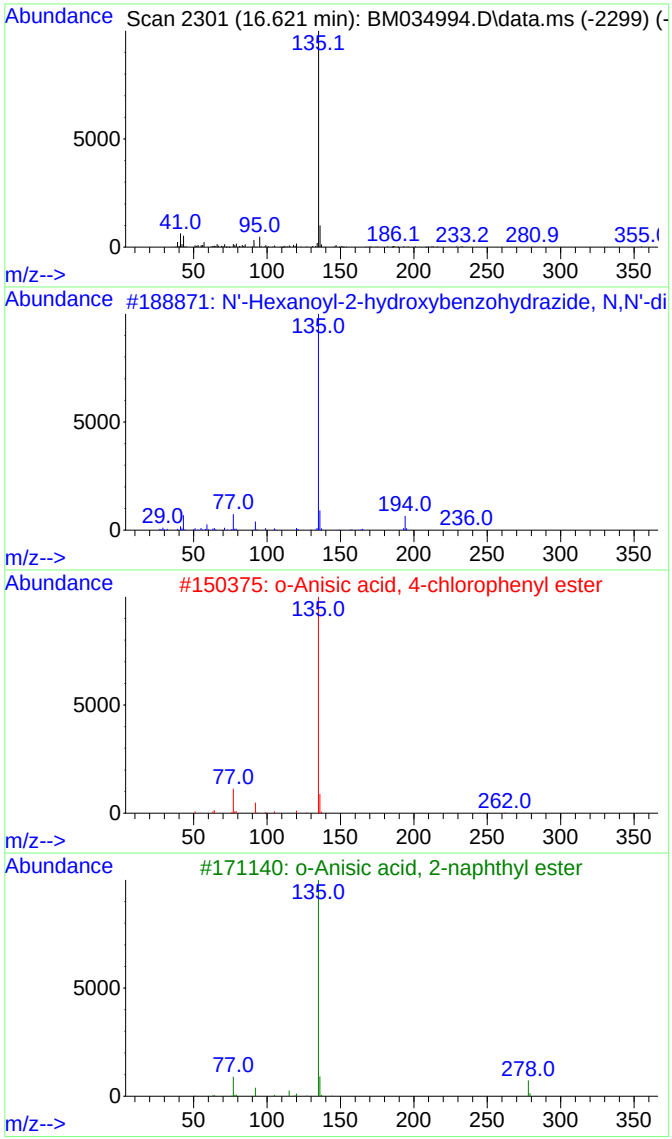
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

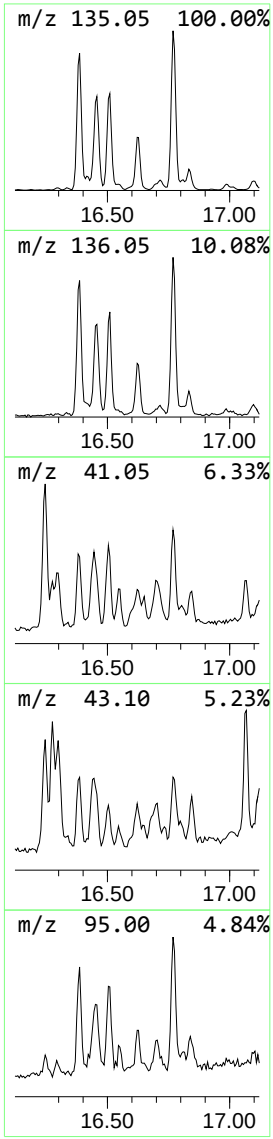
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 N'-Hexanoyl-2-hydroxybenzoh... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.621	2.77 ng/ul	264217	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N'-Hexanoyl-2-hydroxybenzohydraz...	292	C16H24N2O3	1000509-71-7	64
2	o-Anisic acid, 4-chlorophenyl ester	262	C14H11ClO3	1000307-79-0	64
3	o-Anisic acid, 2-naphthyl ester	278	C18H14O3	1000307-79-6	64
4	o-Anisic acid, 4-cyanophenyl ester	253	C15H11NO3	1000307-79-3	50
5	(S)-(7-Bromo-3-isopropyl-1,1-dio...	453	C19H20BrNO5S	1000458-44-1	50





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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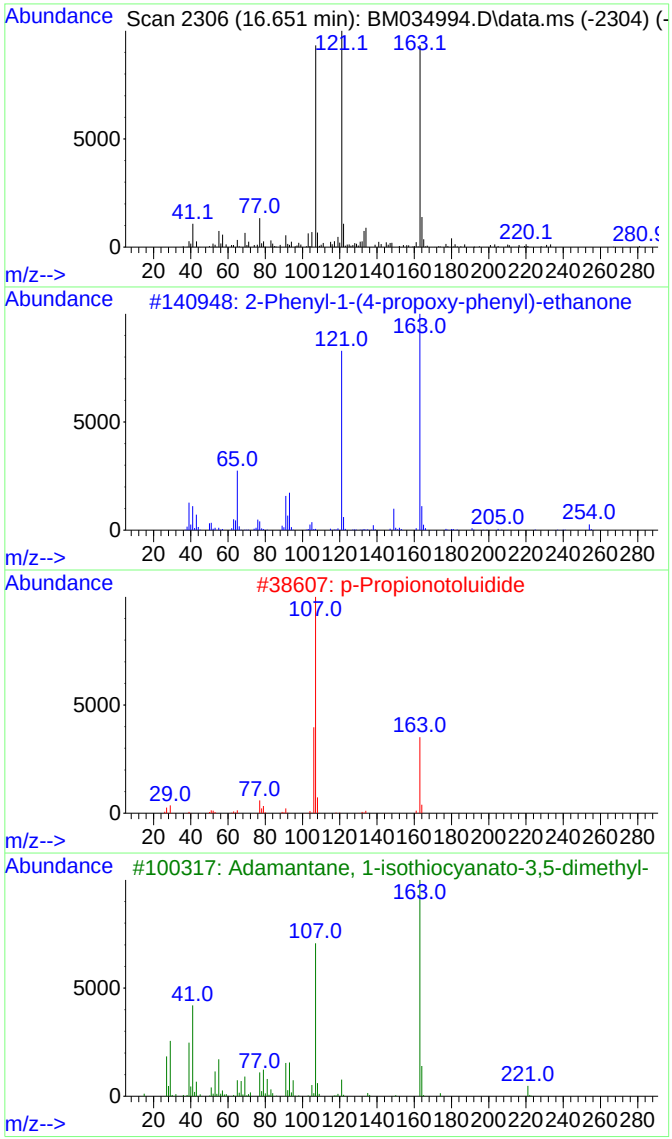
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

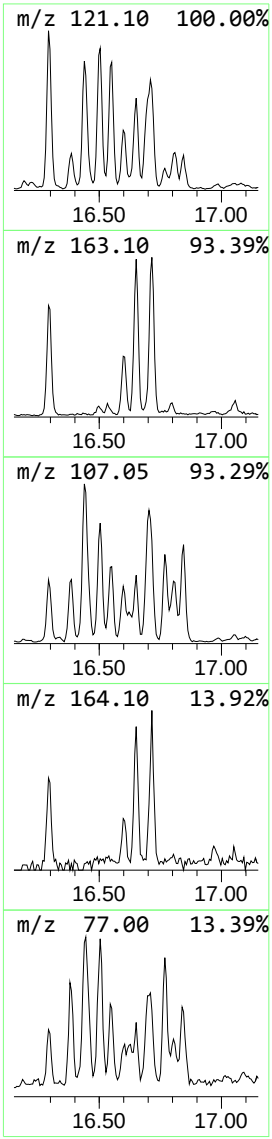
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2-Phenyl-1-(4-propoxy-pheny... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.651	2.22 ng/ul	211950	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenyl-1-(4-propoxy-phenyl)-et...	254	C17H18O2	063192-18-7	50
2	p-Propionotoluidide	163	C10H13NO	002759-55-9	47
3	Adamantane, 1-isothiocyanato-3,5...	221	C13H19NS	136860-49-6	47
4	2-Propanone, 1-(4-methoxyphenyl)-	164	C10H12O2	000122-84-9	25
5	2-Propanone, 1-(4-methoxyphenyl)-	164	C10H12O2	000122-84-9	25





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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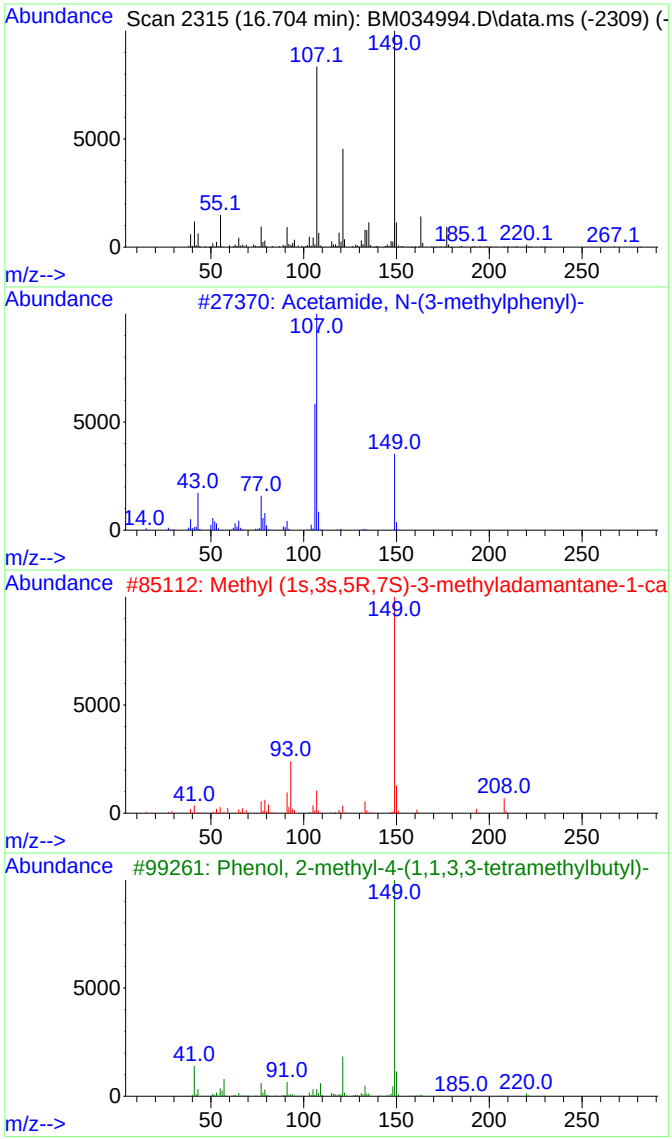
Quant Title : SVOA CALIBRATION

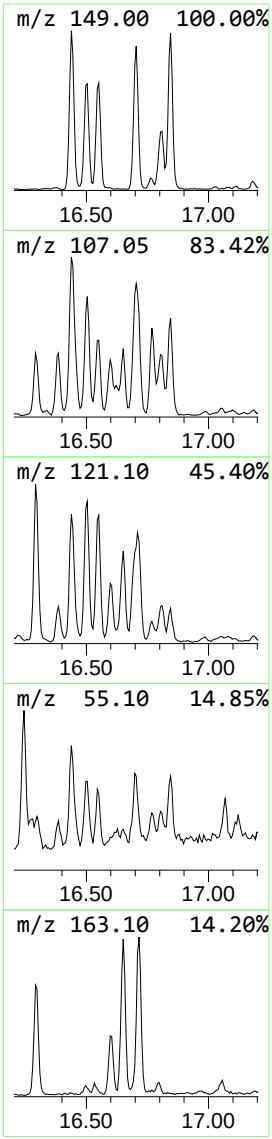
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 unknown-04 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.704	7.97 ng/ul	761063	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
2	Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	38
3	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	38
4	3,4-Methylenedioxypropiophenone	178	C10H10O3	028281-49-4	38
5	Benzene, 1-(1,1-dimethylethyl)-4...	164	C11H16O	005396-38-3	35





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Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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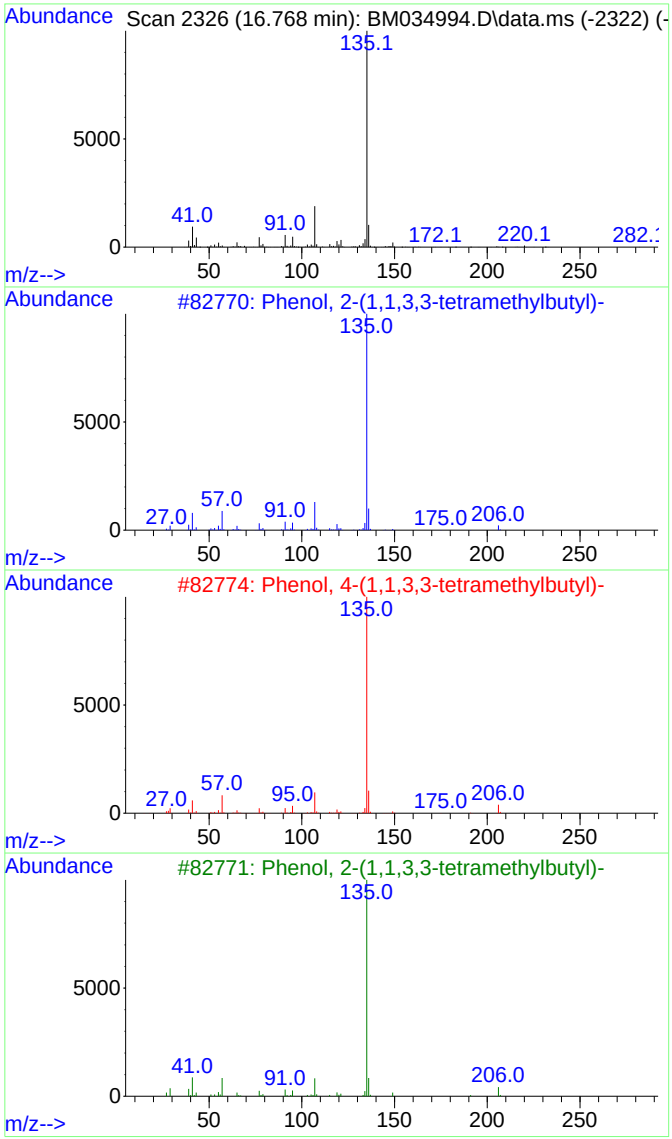
Quant Title : SVOA CALIBRATION

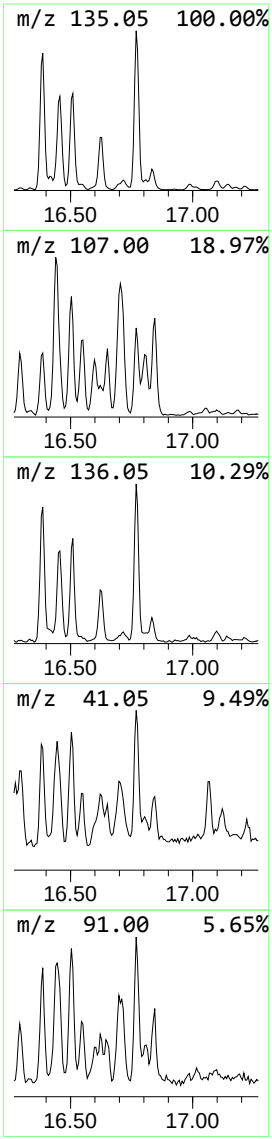
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.768	7.19 ng/ul	686371	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	86
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
3	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
4	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
5	Hexestrol	270	C18H22O2	000084-16-2	72





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ALS Vial : 19 Sample Multiplier: 1

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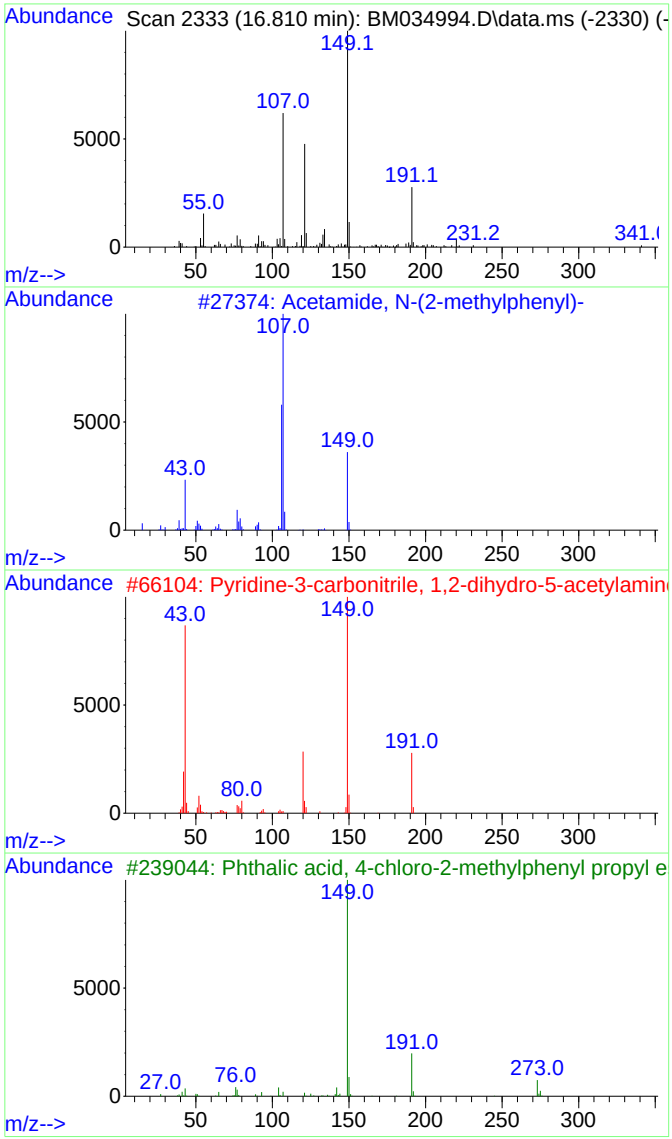
Quant Title : SVOA CALIBRATION

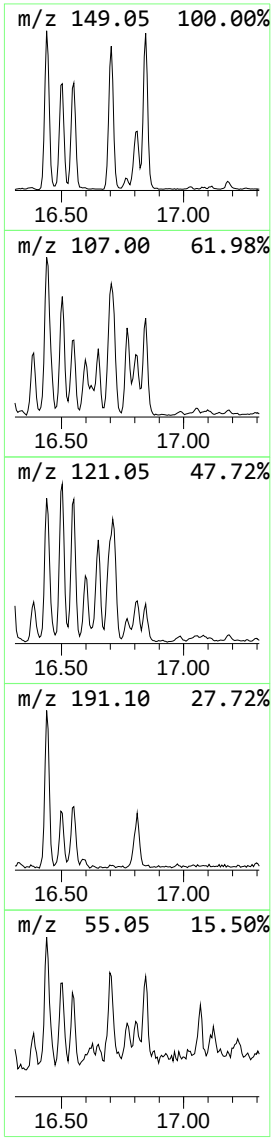
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Acetamide, N-(2-methylphenyl)- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.810	2.42 ng/ul	230754	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	58
2	Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	52
3	Phthalic acid, 4-chloro-2-methyl...	332	C18H17ClO4	1010356-65-9	50
4	Phthalic acid, 2-naphthyl propyl...	334	C21H18O4	1000315-23-9	47
5	Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	47





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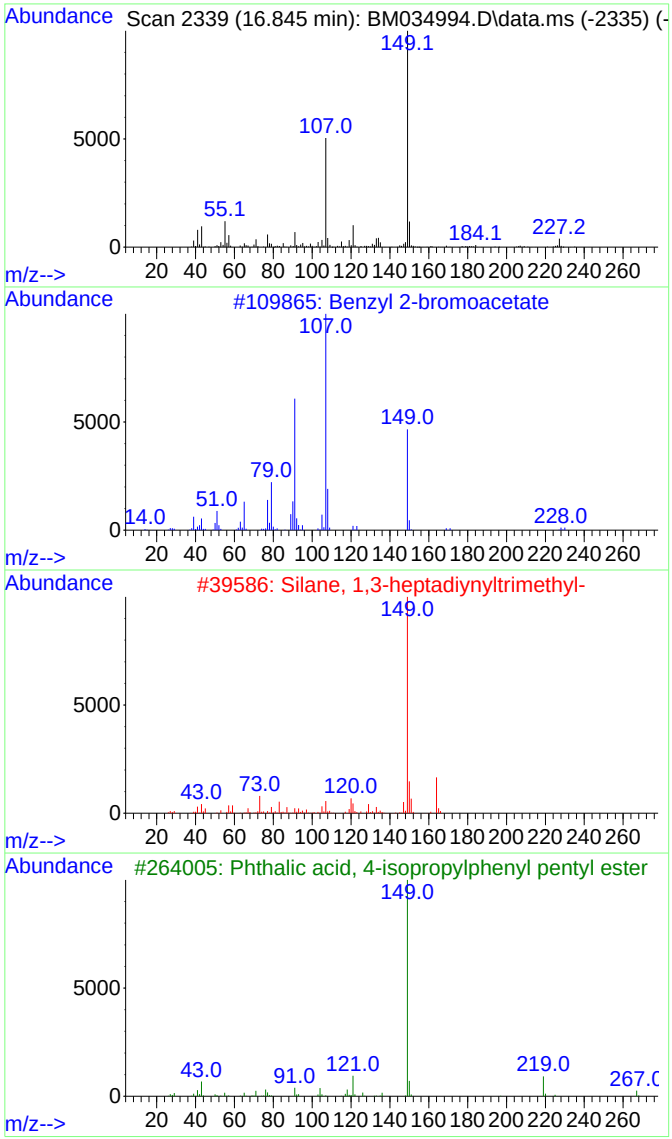
Quant Title : SVOA CALIBRATION

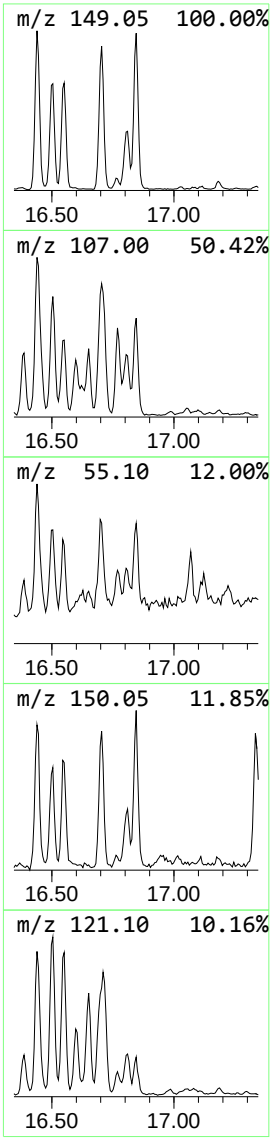
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzyl 2-bromoacetate Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.845	4.72 ng/ul	450906	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzyl 2-bromoacetate	228	C9H9BrO2	005437-45-6	64
2	Silane, 1,3-heptadiynyltrimethyl-	164	C10H16Si	019024-47-6	53
3	Phthalic acid, 4-isopropylphenyl...	354	C22H26O4	1000315-22-2	47
4	7-Methylthieno[3,2-b]pyridine	149	C8H7NS	013362-83-9	47
5	Phthalic acid, di(2-chloropropyl...	318	C14H16Cl2O4	1000356-84-0	47





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Data File : BM034994.D
Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

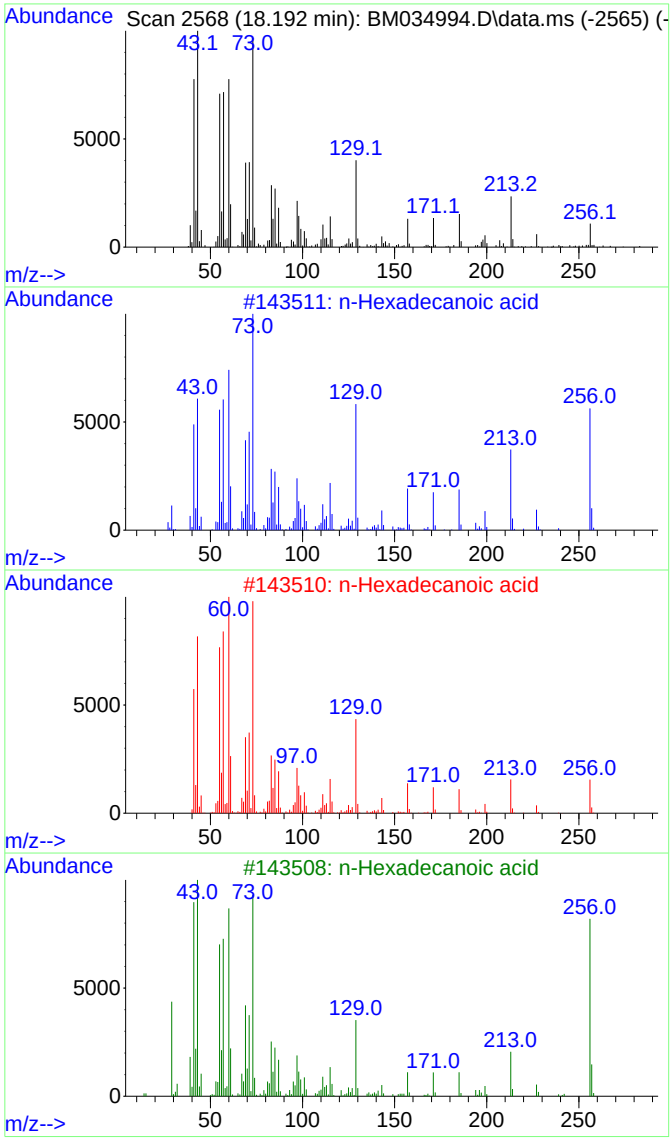
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

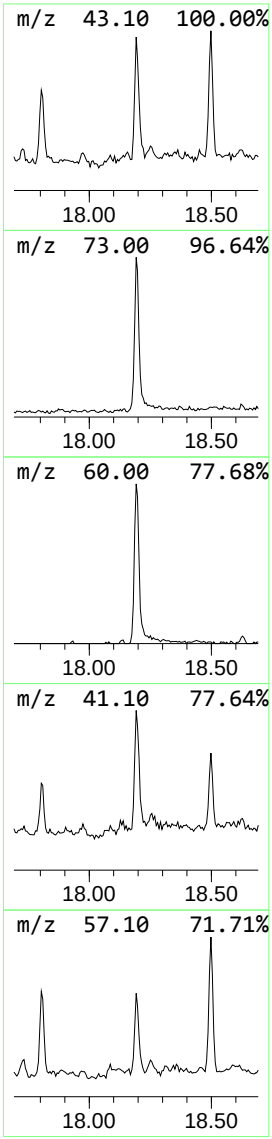
****

Peak Number 17 n-Hexadecanoic acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	4.04 ng/ul	385811	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	87





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Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

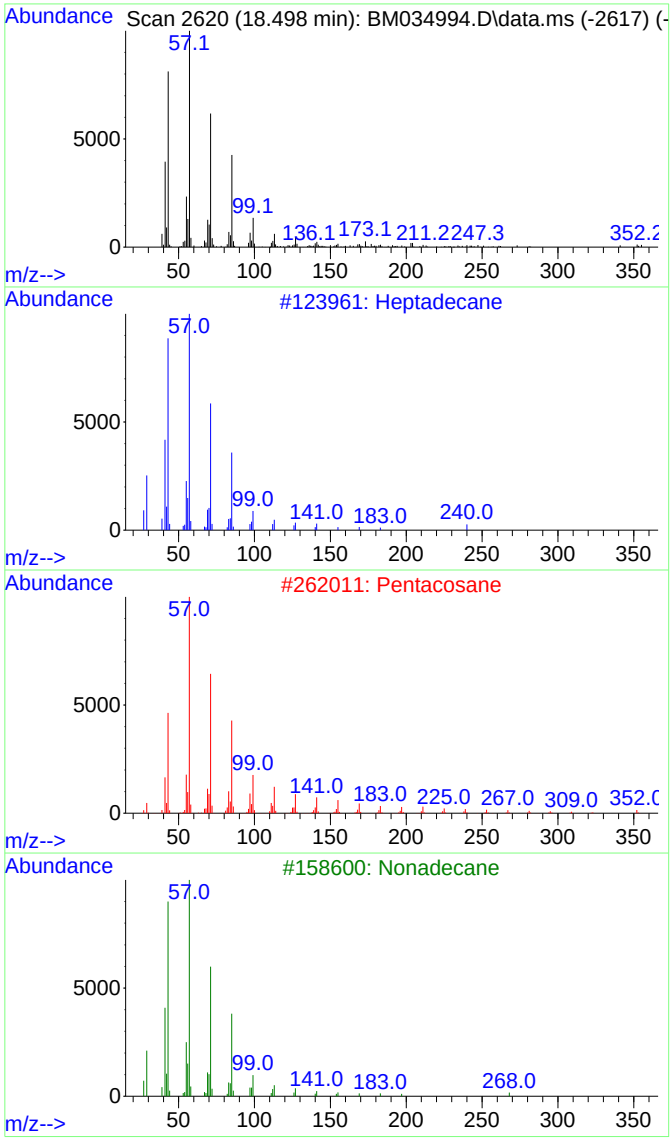
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

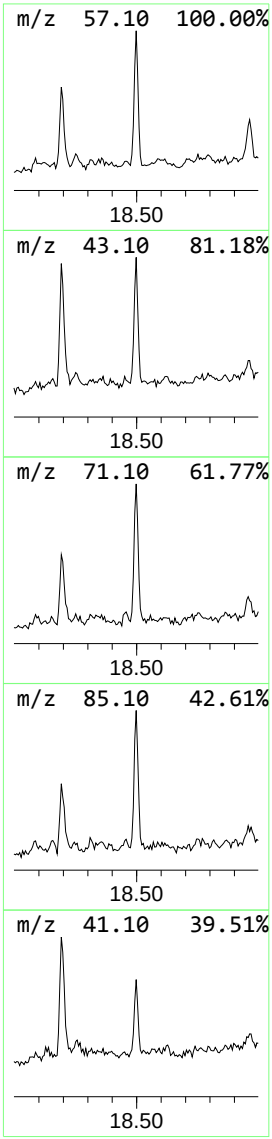
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.498	2.09 ng/ul	199827	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Pentacosane	352	C25H52	000629-99-2	93
3	Nonadecane	268	C19H40	000629-92-5	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Heptacosane	380	C27H56	000593-49-7	91





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

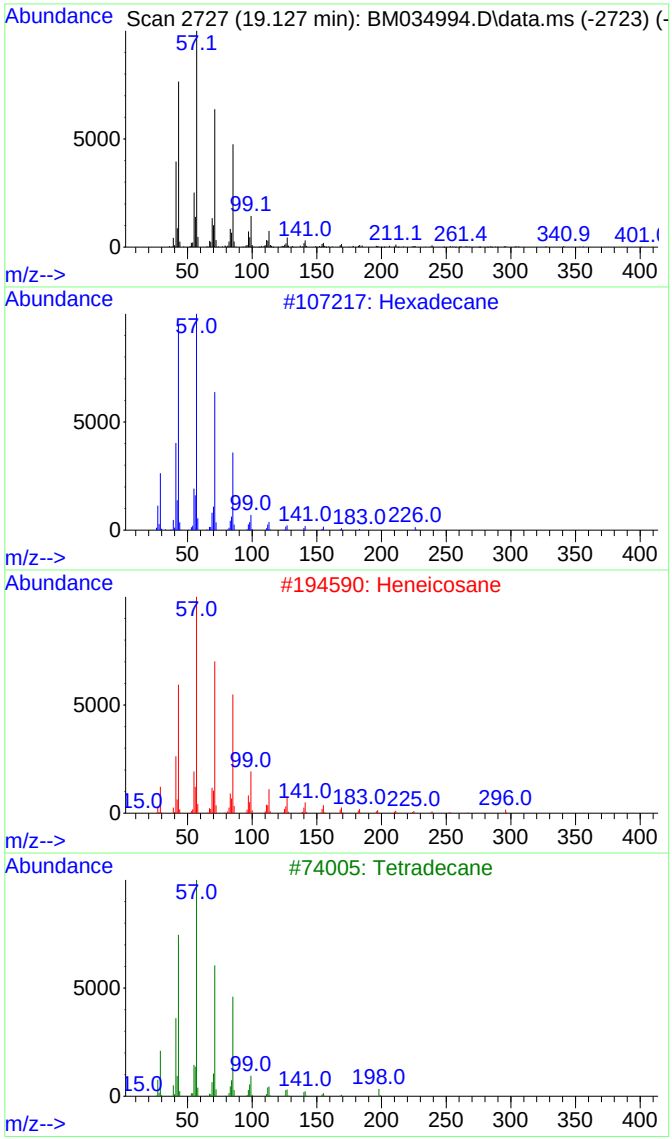
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

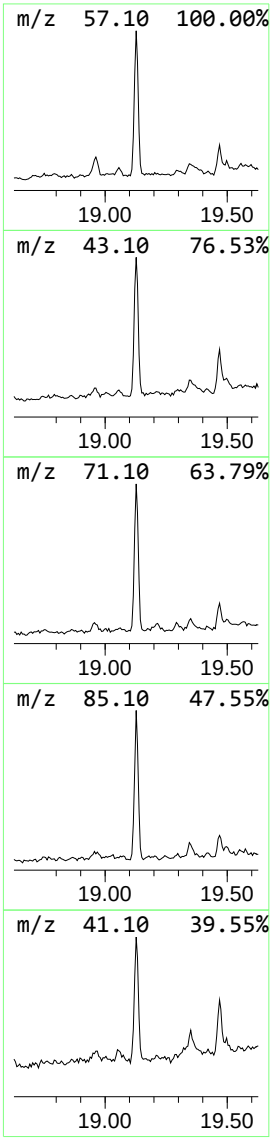
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.127	4.49 ng/ul	428884	Phenanthrene-d10	17.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	96
2	Heneicosane	296	C21H44	000629-94-7	96
3	Tetradecane	198	C14H30	000629-59-4	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Nonadecane	268	C19H40	000629-92-5	91





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

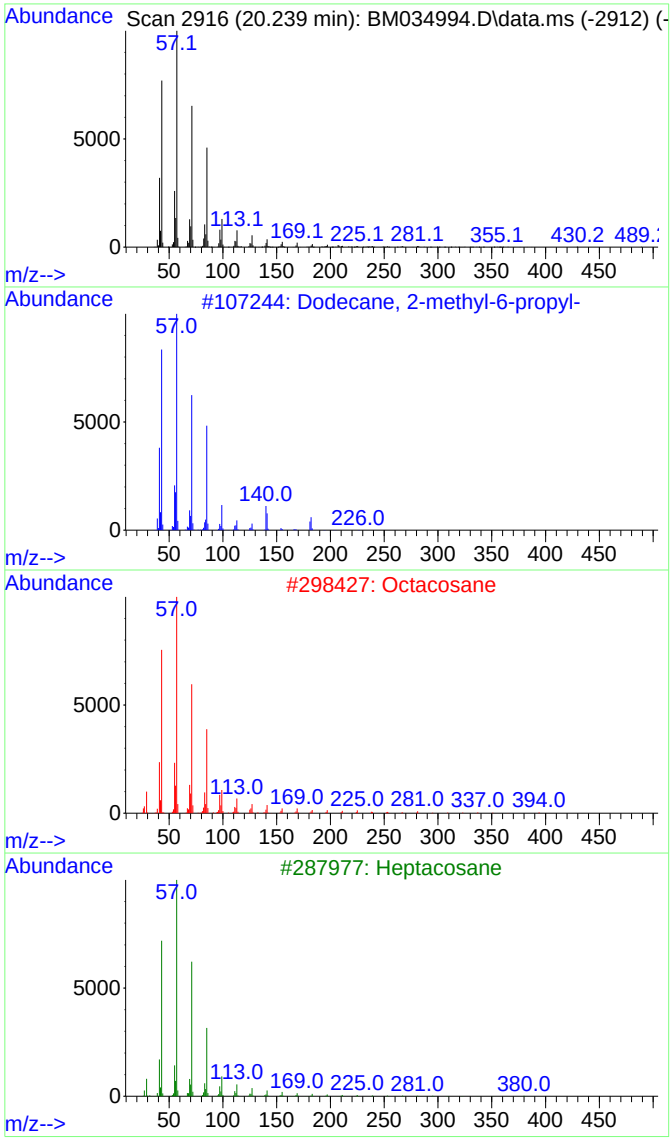
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

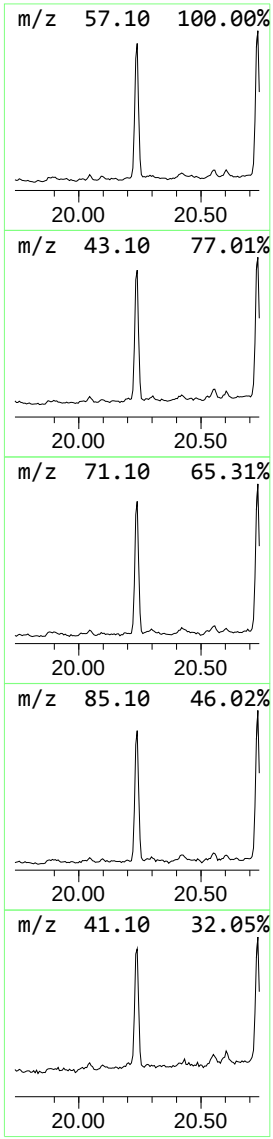
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.239	7.92 ng/ul	856938	Chrysene-d12	21.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
2	Octacosane	394	C28H58	000630-02-4	91
3	Heptacosane	380	C27H56	000593-49-7	91
4	Eicosane	282	C20H42	000112-95-8	91
5	Heneicosane	296	C21H44	000629-94-7	91





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

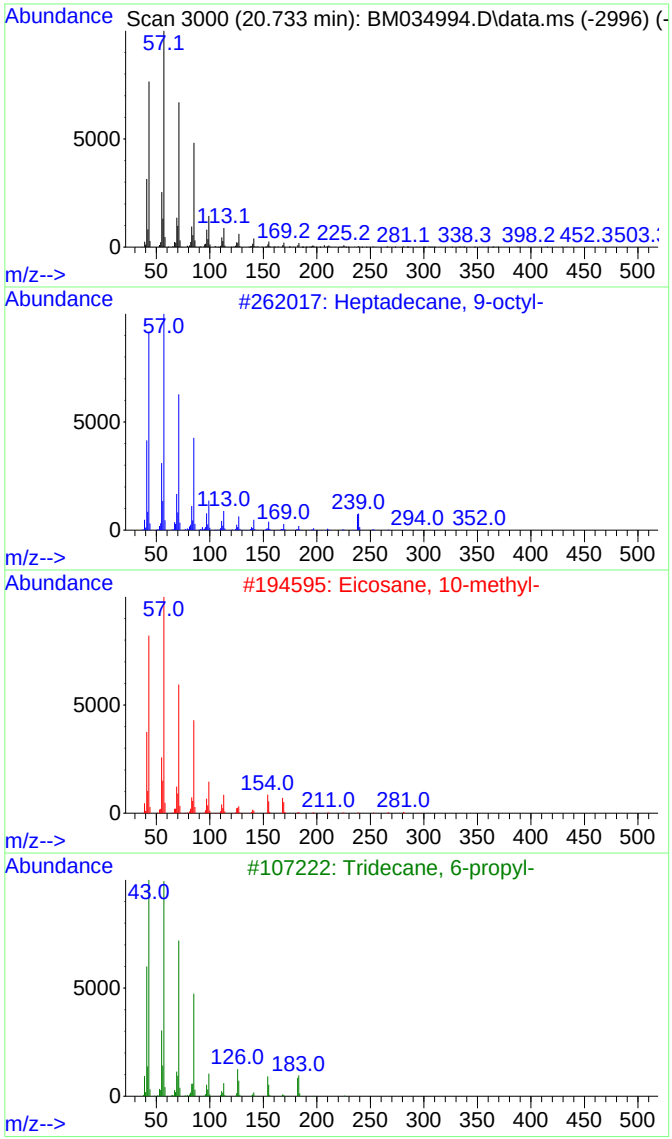
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

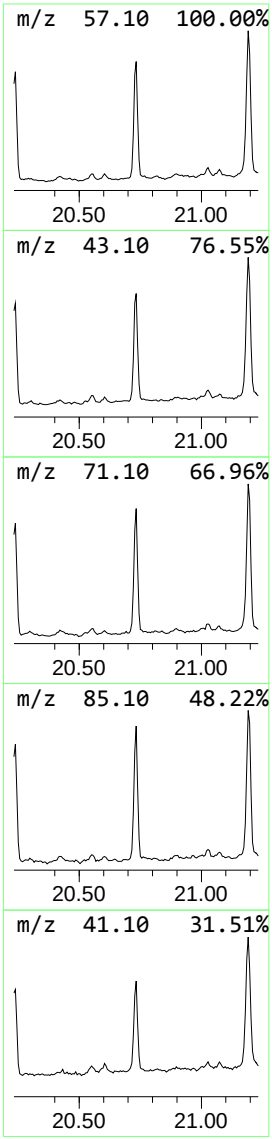
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.733	8.37 ng/ul	904899	Chrysene-d12	21.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3	Tridecane, 6-propyl-	226	C16H34	055045-10-8	93
4	Heneicosane	296	C21H44	000629-94-7	91
5	Hexacosane	366	C26H54	000630-01-3	91





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

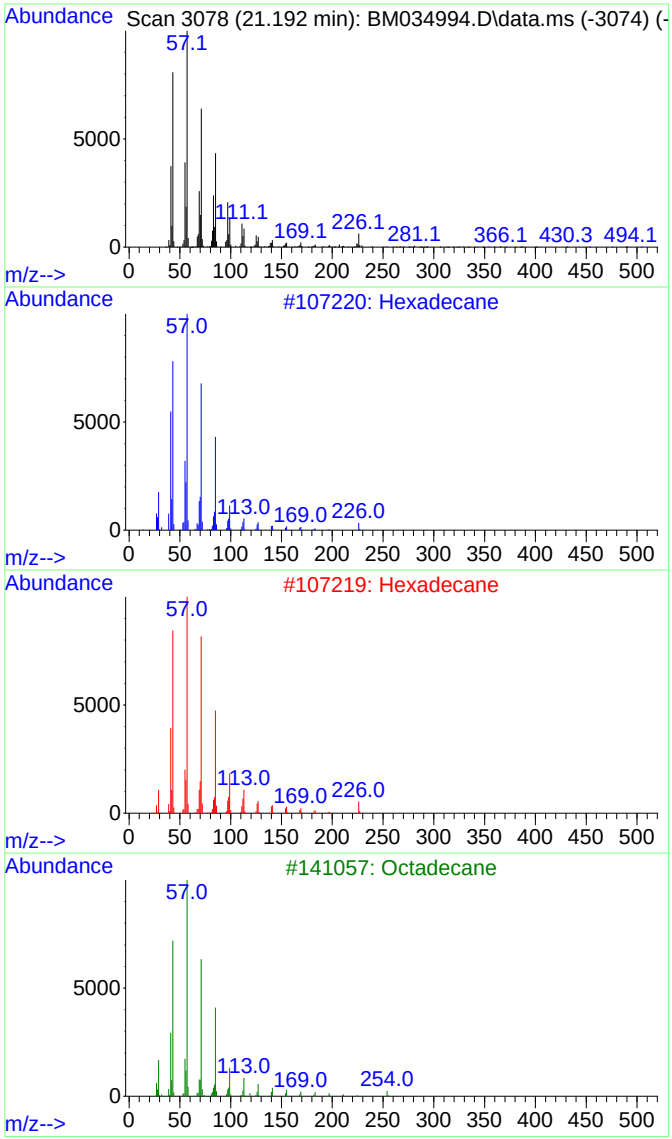
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

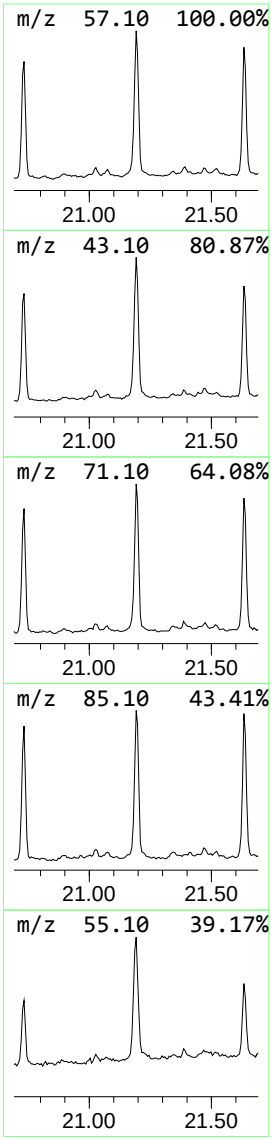
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.192	14.77 ng/ul	1597170	Chrysene-d12	21.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	96
2	Hexadecane	226	C16H34	000544-76-3	93
3	Octadecane	254	C18H38	000593-45-3	93
4	Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	91
5	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91





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Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

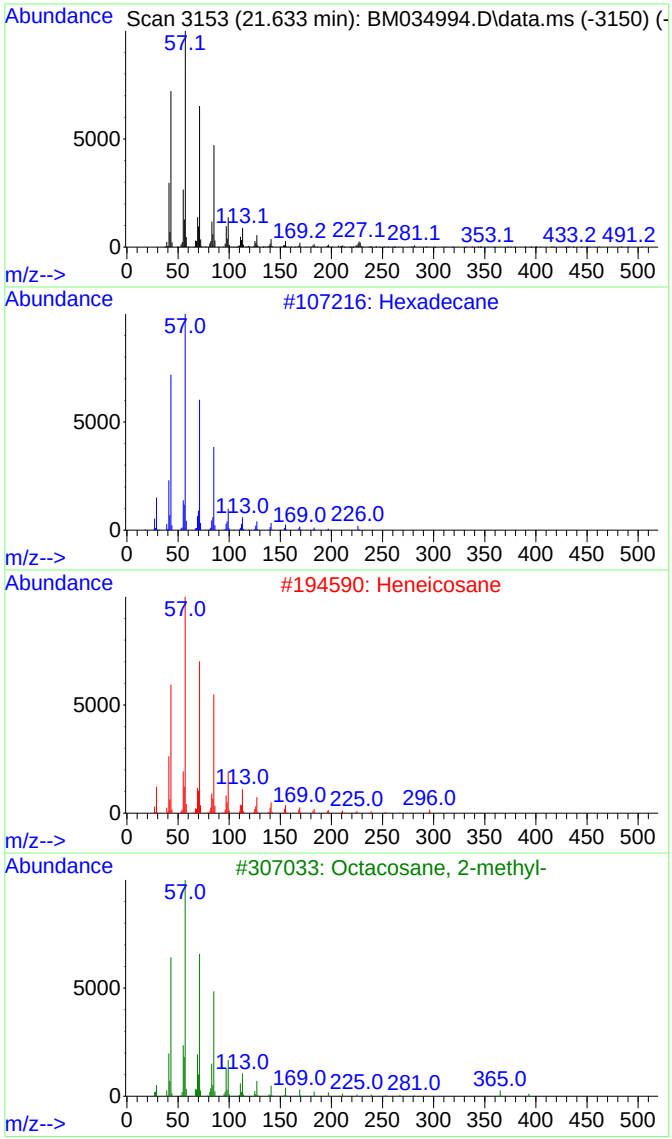
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

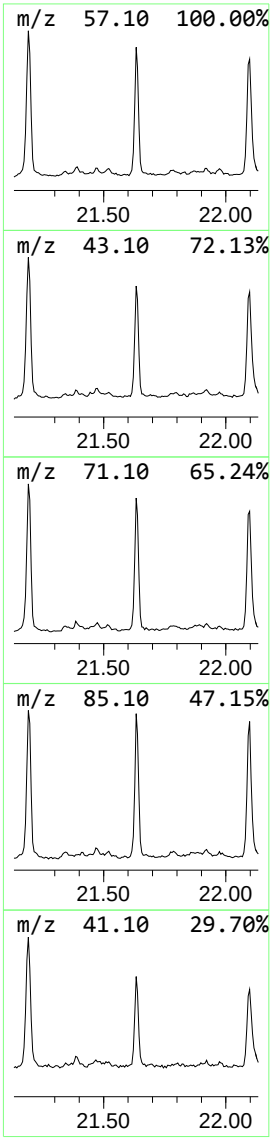
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.633	9.18 ng/ul	993412	Chrysene-d12	21.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	96
2	Heneicosane	296	C21H44	000629-94-7	91
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4	Heptacosane	380	C27H56	000593-49-7	91
5	Octacosane	394	C28H58	000630-02-4	91





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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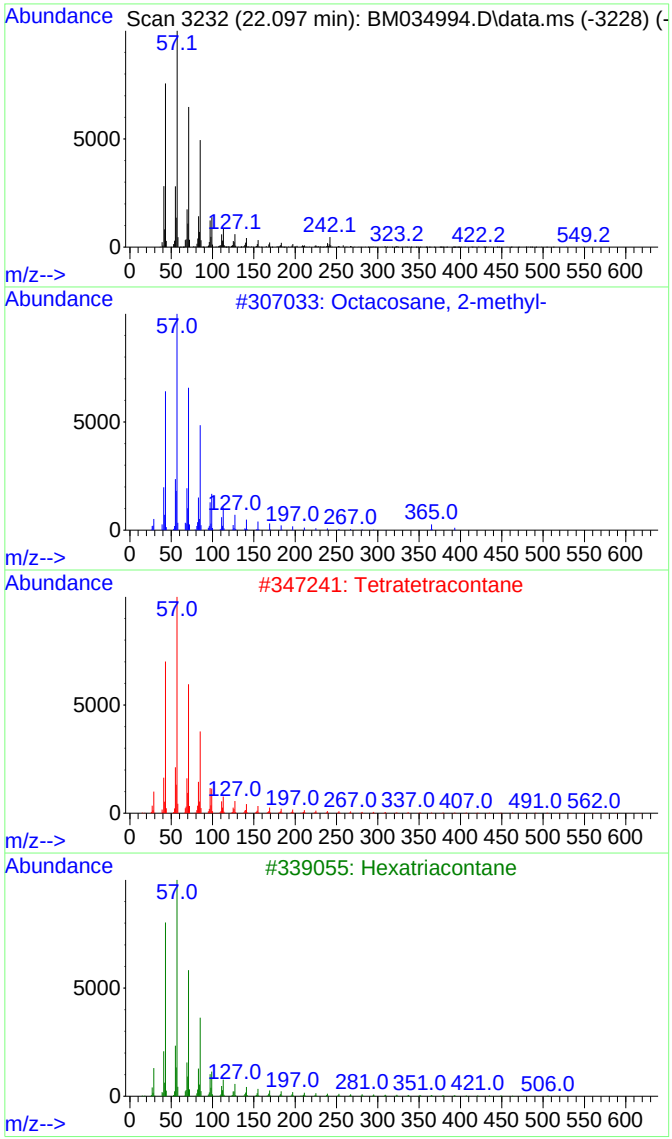
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

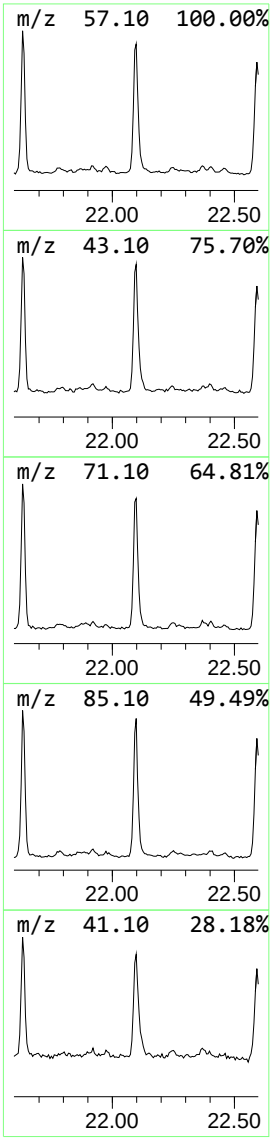
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.097	12.93 ng/ul	1398040	Chrysene-d12	21.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Hexatriacontane	507	C36H74	000630-06-8	91
4	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	91





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Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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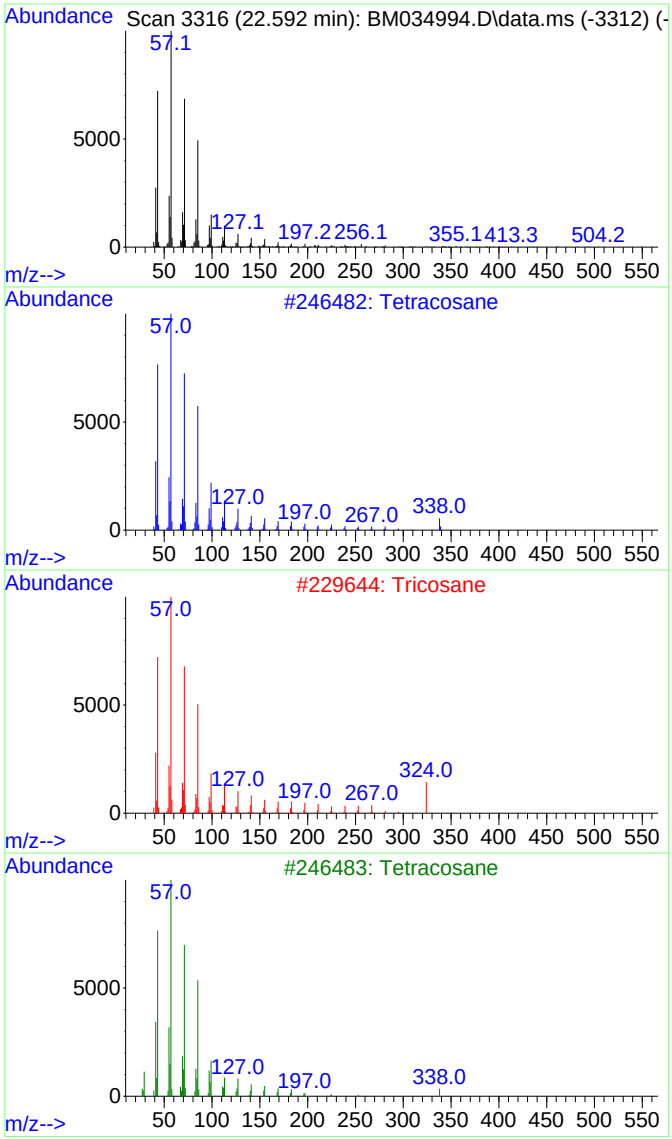
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

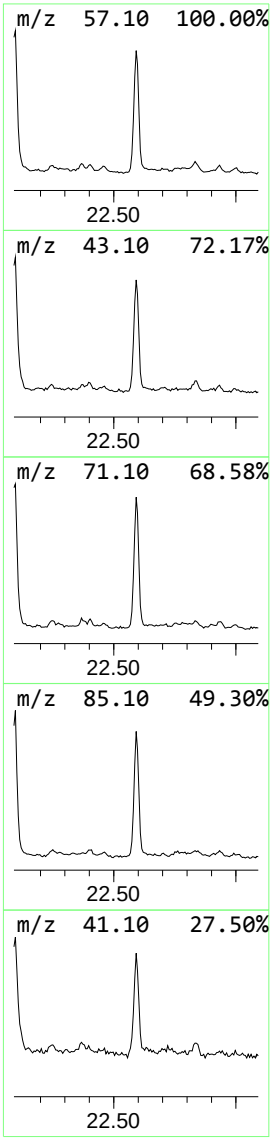
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.592	9.51 ng/ul	1028160	Chrysene-d12	21.468

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	96
2	Tricosane	324	C23H48	000638-67-5	95
3	Tetracosane	338	C24H50	000646-31-1	95
4	Heptadecane	240	C17H36	000629-78-7	95
5	Eicosane, 9-octyl-	394	C28H58	013475-77-9	94





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Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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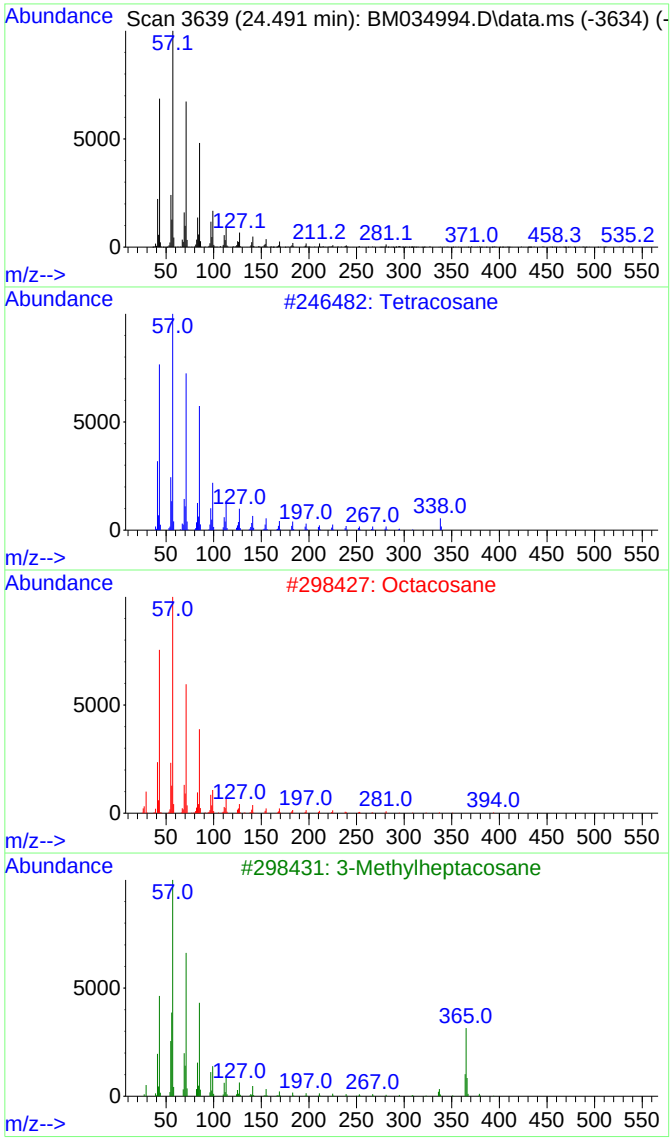
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

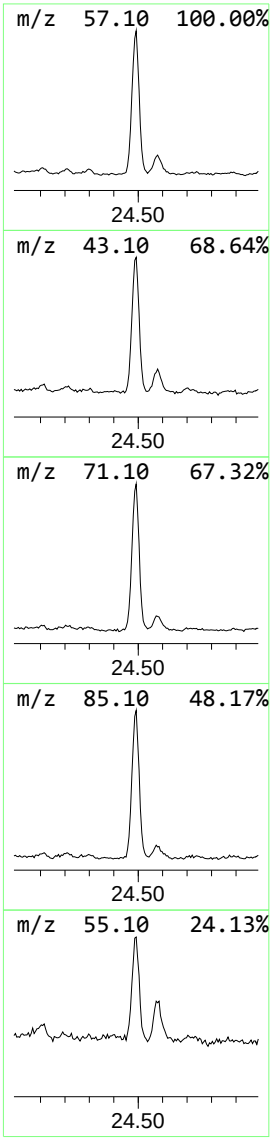
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.491	16.40 ng/ul	1606330	Perylene-d12	23.856

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Octacosane	394	C28H58	000630-02-4	95
3	3-Methylheptacosane	394	C28H58	014167-66-9	93
4	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93





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Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

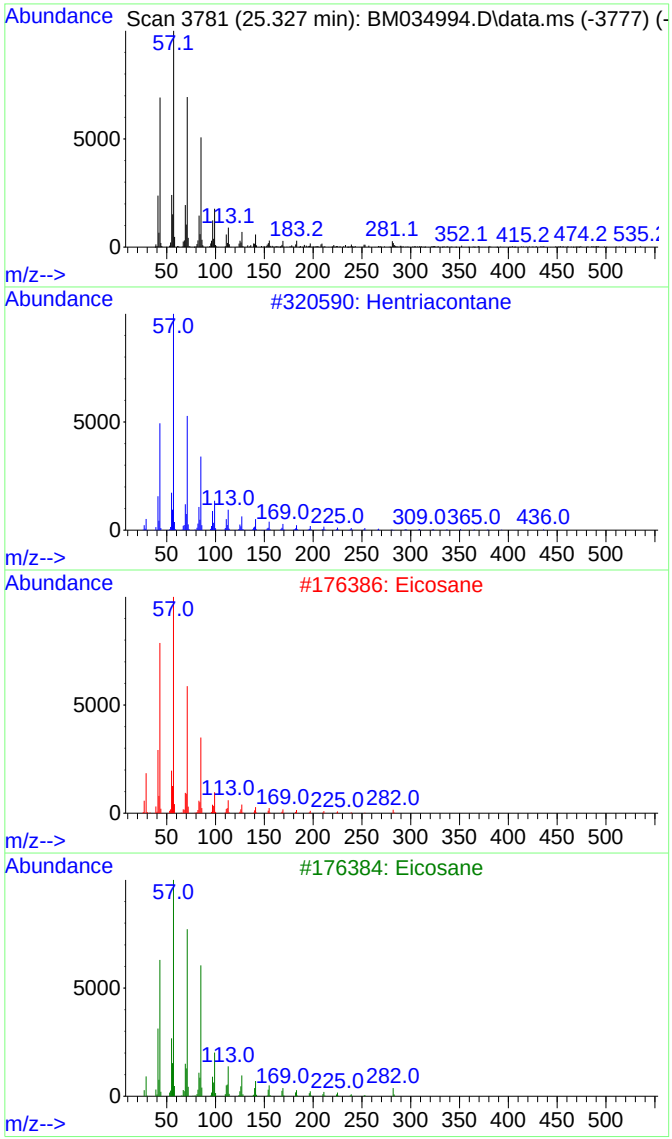
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION

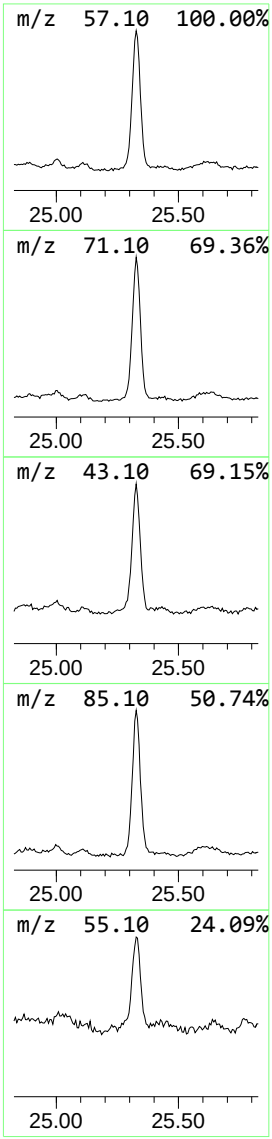
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.327	11.24 ng/ul	1100680	Perylene-d12	23.856

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	96
2	Eicosane	282	C20H42	000112-95-8	95
3	Eicosane	282	C20H42	000112-95-8	93
4	Octacosane	394	C28H58	000630-02-4	90
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM034994.D
Acq On : 11 May 2022 20:33
Operator : CG/JU
Sample : N2603-17
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW497

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.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
Methacrylamide	3.887	2.3	ng/ul	141277	1	7.922	1243480	20.0
Benzyl alcohol	8.157	6.0	ng/ul	370980	1	7.922	1243480	20.0
Ethanol, 2-(2-b...	10.504	2.3	ng/ul	199226	2	10.722	1700780	20.0
unknown-01	16.245	6.8	ng/ul	645947	4	17.292	1909160	20.0
Acetamide, N-(2...	16.298	5.1	ng/ul	484709	4	17.292	1909160	20.0
Phenol, 2-(1,1,...	16.386	5.8	ng/ul	557163	4	17.292	1909160	20.0
unknown-02	16.445	10.5	ng/ul	999901	4	17.292	1909160	20.0
4-(7-Methylocty...	16.504	8.0	ng/ul	761886	4	17.292	1909160	20.0
Hexestrol dimet...	16.551	4.2	ng/ul	401486	4	17.292	1909160	20.0
unknown-03	16.604	2.1	ng/ul	200374	4	17.292	1909160	20.0
N'-Hexanoyl-2-h...	16.621	2.8	ng/ul	264217	4	17.292	1909160	20.0
2-Phenyl-1-(4-p...	16.651	2.2	ng/ul	211950	4	17.292	1909160	20.0
unknown-04	16.704	8.0	ng/ul	761063	4	17.292	1909160	20.0
Phenol, 4-(1,1,...	16.768	7.2	ng/ul	686371	4	17.292	1909160	20.0
Acetamide, N-(2...	16.810	2.4	ng/ul	230754	4	17.292	1909160	20.0
Benzyl 2-bromoa...	16.845	4.7	ng/ul	450906	4	17.292	1909160	20.0
n-Hexadecanoic ...	18.192	4.0	ng/ul	385811	4	17.292	1909160	20.0
(DEL) Alkane: S...	18.498	2.1	ng/ul	199827	4	17.292	1909160	20.0
(DEL) Alkane: S...	19.127	4.5	ng/ul	428884	4	17.292	1909160	20.0
(DEL) Alkane: S...	20.239	7.9	ng/ul	856938	5	21.468	2163240	20.0
(DEL) Alkane: S...	20.733	8.4	ng/ul	904899	5	21.468	2163240	20.0
(DEL) Alkane: S...	21.192	14.8	ng/ul	1597170	5	21.468	2163240	20.0
(DEL) Alkane: S...	21.633	9.2	ng/ul	993412	5	21.468	2163240	20.0
(DEL) Alkane: S...	22.097	12.9	ng/ul	1398040	5	21.468	2163240	20.0
(DEL) Alkane: S...	22.592	9.5	ng/ul	1028160	5	21.468	2163240	20.0
(DEL) Alkane: S...	24.491	16.4	ng/ul	1606330	6	23.856	1959330	20.0
(DEL) Alkane: S...	25.327	11.2	ng/ul	1100680	6	23.856	1959330	20.0