LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

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-BP072421.M
Title : SVOA CALIBRATION

Signal : TIC: BP006608.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.276	29	32	41	rVB2	61190	97976	1.71%	0.108%
2	3.681	97	101	115	rBV	678135	1122627	19.63%	1.243%
3	3.781	115	118	126	rVB	92744	136132	2.38%	0.151%
4	6.934	647	654	665	rVB	1052271	1740796	30.44%	1.927%
5	7.099	676	682	695	rVB	883767	1420979	24.85%	1.573%
6	7.287	707	714	722	rBV	1131540	1753477	30.66%	1.941%
7	7.752	787	793	800	rVB	615355	957322	16.74%	1.060%
8	8.463	907	914	922	rBV	872318	1402401	24.52%	1.553%
9	8.905	982	989	997	rBV	1098497	1815777	31.75%	2.010%
10	9.622	1104	1111	1126	rBV	943628	1544934	27.01%	1.710%
11	10.158	1195	1202	1212	rBV	1257808	2054439	35.92%	2.275%
12	10.252	1212	1218	1224	rVV2	70050	121493	2.12%	0.135%
13	10.328	1224	1231	1242	rVB	503604	844300	14.76%	0.935%
14	10.534	1258	1266	1272	rBV	849866	1428672	24.98%	1.582%
15	10.581	1272	1274	1280	rVB	41132	58940	1.03%	0.065%
16	10.681	1283	1291	1297	rBV	893830	1631259	28.52%	1.806%
17	11.963	1503	1509	1515	rVV	36204	53484	0.94%	0.059%
18	12.122	1530	1536	1541	rVV2	28960	54624	0.96%	0.060%
19	12.199	1541	1549	1556	rVB	46754	80605	1.41%	0.089%
20	13.216	1715	1722	1727	rBV	56319	83726	1.46%	0.093%
21	13.799	1816	1821	1830	rVV	2192990	3074006	53.75%	3.403%
22	14.075	1858	1868	1885	rBV	2498472	3922047	68.58%	4.342%
23	14.293	1900	1905	1910	rVB	49016	66409	1.16%	0.074%
24	14.381	1910	1920	1926	rBV	1137957	1722428	30.12%	1.907%
25	14.593	1950	1956	1976	rVV	828145	1418098	24.80%	1.570%
26	14.746	1976	1982	1985	rBV3	39477	58671	1.03%	0.065%
27	15.140	2043	2049	2053	rBV	132089	178309	3.12%	0.197%
28	15.246	2063	2067	2072	rVB	55676	70270	1.23%	0.078%
29	15.375	2082	2089	2096	rBV	3068857	4467282	78.11%	4.946%
30	15.440	2096	2100	2105	rVV5	33591	57125	1.00%	0.063%
31	15.504	2105	2111	2123	rVB	755746	1025707	17.94%	1.136%
32	15.728	2143	2149	2155	rBV2	25045	54713	0.96%	0.061%
33	16.110	2210	2214	2221	rVB	73849	104933	1.83%	0.116%
34	16.351	2246	2255	2260	rBV2	62597	113038	1.98%	0.125%
35	16.681	2305	2311	2314	rVV4	30083	60225	1.05%	0.067%
36	16.792	2325	2330	2338	rBV2	32505	68797	1.20%	0.076%

LSC Area Percent Report

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Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

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-BP072421.M
Title : SVOA CALIBRATION

37	16.910	2346	2350	2354	rVV2	63486	103398	1.81%	0.114%
38	17.045	2370	2373	2377	rBV	41518	50210	0.88%	0.056%
39	17.134	2380	2388	2392	rBV	1312723	1906994	33.34%	2.111%
40	17.175	2392	2395	2399	rVV	410524	545259	9.53%	0.604%
41	17.234	2399	2405	2416	rVB2	3120193	4617897	80.75%	5.113%
42	17.322	2416	2420	2424	rBV6	31813	55346	0.97%	0.061%
43	17.440	2432	2440	2447	rBV7	51063	131365	2.30%	0.145%
44	17.522	2448	2454	2455	rBV5	40552	61985	1.08%	0.069%
45	17.645	2468	2475	2480	rBV3	86850	161814	2.83%	0.179%
46	17.710	2482	2486	2490	rVV3	104119	139408	2.44%	0.154%
47	17.845	2506	2509	2514	rVB3	184317	247469	4.33%	0.274%
48	17.916	2514	2521	2525	rBV6	91023	188127	3.29%	0.208%
49	17.963	2525	2529	2533	rBV4	88712	189520	3.31%	0.210%
50	18.004	2533	2536	2539	rVV2	167399	223466	3.91%	0.247%
51	18.039	2539	2542	2550	rVB2	331785	507068	8.87%	0.561%
52	18.110	2550	2554	2558	rBV2	229460	307857	5.38%	0.341%
53	18.181	2563	2566	2571	rVV3	157546	255657	4.47%	0.283%
54	18.228	2571	2574	2579	rVB	191794	249744	4.37%	0.277%
55	18.287	2579	2584	2587	rBV5	43425	80257	1.40%	0.089%
56	18.328	2587	2591	2597	rBV3	110189	174452	3.05%	0.193%
57	18.428	2603	2608	2615	rVB4	158339	305878	5.35%	0.339%
58	18.492	2615	2619	2623	rVV	116377	162736	2.85%	0.180%
59	18.534	2623	2626	2634	rVB3	68114	111014	1.94%	0.123%
60	18.816	2669	2674	2679	rVB4	77761	143969	2.52%	0.159%
61	18.898	2684	2688	2692	rVV	61975	94819	1.66%	0.105%
62	18.939	2692	2695	2700	rVB3	73713	104182	1.82%	0.115%
63	19.051	2707	2714	2718	rBV2	61320	146937	2.57%	0.163%
64	19.151	2722	2731	2739	rBV	1255452	1704106	29.80%	1.887%
65	19.275	2750	2752	2760	rVB3	52806	97940	1.71%	0.108%
66	19.481	2780	2787	2790	rBV	3866267	5719023	100.00%	6.332%
67	19.575	2800	2803	2809	rVB2	65234	101209	1.77%	0.112%
68	19.681	2818	2821	2826	rVB	51385	65386	1.14%	0.072%
69	19.822	2840	2845	2851	rBV6	75103	196334	3.43%	0.217%
70	19.986	2870	2873	2876	rBV2	113517	135685	2.37%	0.150%
71	20.022	2876	2879	2882	rVB	104481	117520	2.05%	0.130%
72	20.086	2887	2890	2893	rBV	73627	87123	1.52%	0.096%
73	20.281	2920	2923	2927	rBV2	93047	107563	1.88%	0.119%
74	20.392	2939	2942	2946	rBV	136707	147414	2.58%	0.163%
75	20.492	2955	2959	2965	rBV2	214201	372619	6.52%	0.413%
76	20.604	2975	2978	2982	rVB	371760	395131	6.91%	0.437%

LSC Area Percent Report

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Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

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_P\METHODS\SFAM-EPA-BP072421.M
Title : SVOA CALIBRATION

77	20.722	2994	2998	3005	rVB2	299664	522207	9.13%	0.578%
78	20.839	3015	3018	3022	rBV3	192465	213032	3.72%	0.236%
79	20.922	3030	3032	3034	rBV	79021	74720	1.31%	0.083%
80	20.963	3035	3039	3043	rVV2	619239	794700	13.90%	0.880%
81	21.128	3064	3067	3069	rBV3	141995	184078	3.22%	0.204%
82	21.157	3069	3072	3075	rVV2	225137	264965	4.63%	0.293%
83	21.228	3076	3084	3088	rVV2	1767421	3318169	58.02%	3.674%
84	21.269	3088	3091	3094	rVV	572436	659451	11.53%	0.730%
85	21.304	3094	3097	3103	rVV5	178908	304656	5.33%	0.337%
86	21.392	3108	3112	3114	rBV2	364600	455963	7.97%	0.505%
87	21.422	3114	3117	3121	rVV2	369357	460127	8.05%	0.509%
88	21.622	3147	3151	3154	rBV2	408608	520101	9.09%	0.576%
89	21.769	3172	3176	3179	rBV	828140	1036962	18.13%	1.148%
90	22.292	3262	3265	3267	rBV2	311814	420756	7.36%	0.466%
91	22.327	3268	3271	3277	rVB	485873	813372	14.22%	0.901%
92	22.863	3357	3362	3374	rVB2	1065660	2572924	44.99%	2.849%
93	23.416	3450	3456	3461	rBV2	2531953	4506234	78.79%	4.989%
94	23.469	3461	3465	3472	rVB3	890904	1763393	30.83%	1.952%
95	23.563	3475	3481	3487	rBV2	1010863	1885350	32.97%	2.087%
96	24.174	3580	3585	3596	rVB	828844	1792307	31.34%	1.984%
97	24.992	3718	3724	3733	rVB	952709	2172711	37.99%	2.406%
98	25.951	3880	3887	3902	rVB	1539700	4468462	78.13%	4.947%
99	27.092	4074	4081	4095	rVB2	985739	3186734	55.72%	3.528%
100	28.433	4303	4309	4322	rVB	743756	2550225	44.59%	2.823%

Sum of corrected areas: 90321501

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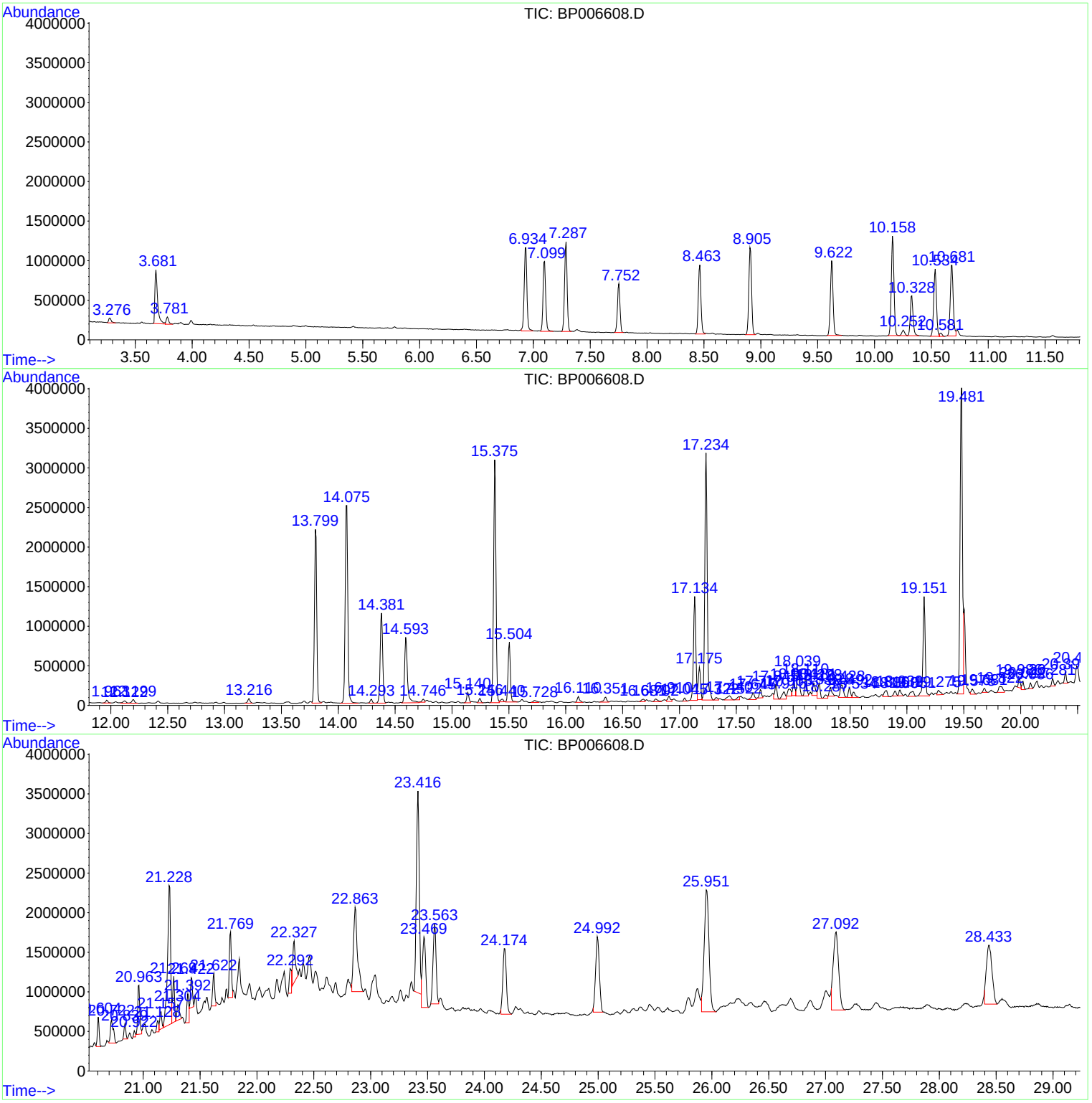
Instrument :
BNA_P
ClientSampleId :
C00A2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006608.D
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Instrument : BNA_P
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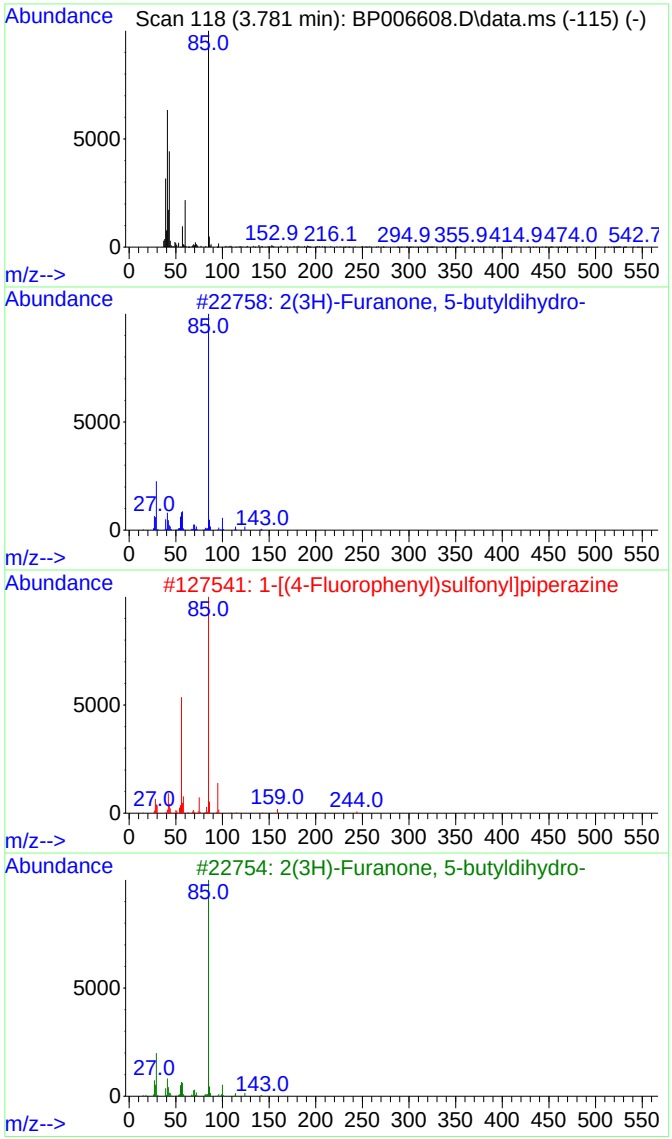
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

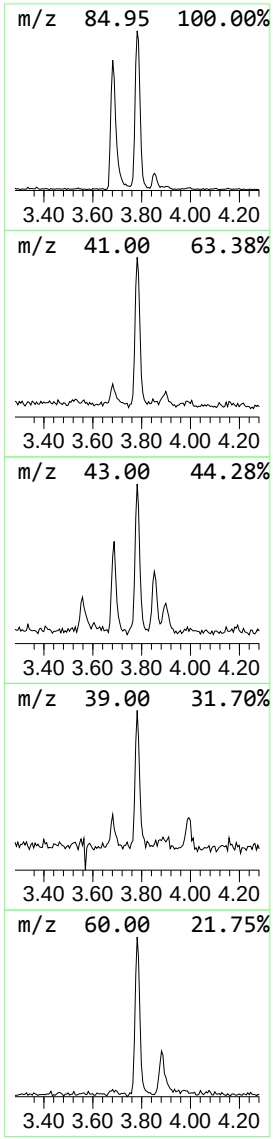
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2(3H)-Furanone, 5-butyldihy... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.780	2.84 ng/ul	136132	1,4-Dichlorobenzene-d4	7.752

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Furanone, 5-butyldihydro-	142	C8H14O2	000104-50-7	53
2	1-[(4-Fluorophenyl)sulfonyl]pipe...	244	C10H13FN2O2S	1000486-20-7	42
3	2(3H)-Furanone, 5-butyldihydro-	142	C8H14O2	000104-50-7	42
4	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
5	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38





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Instrument :
BNA_P
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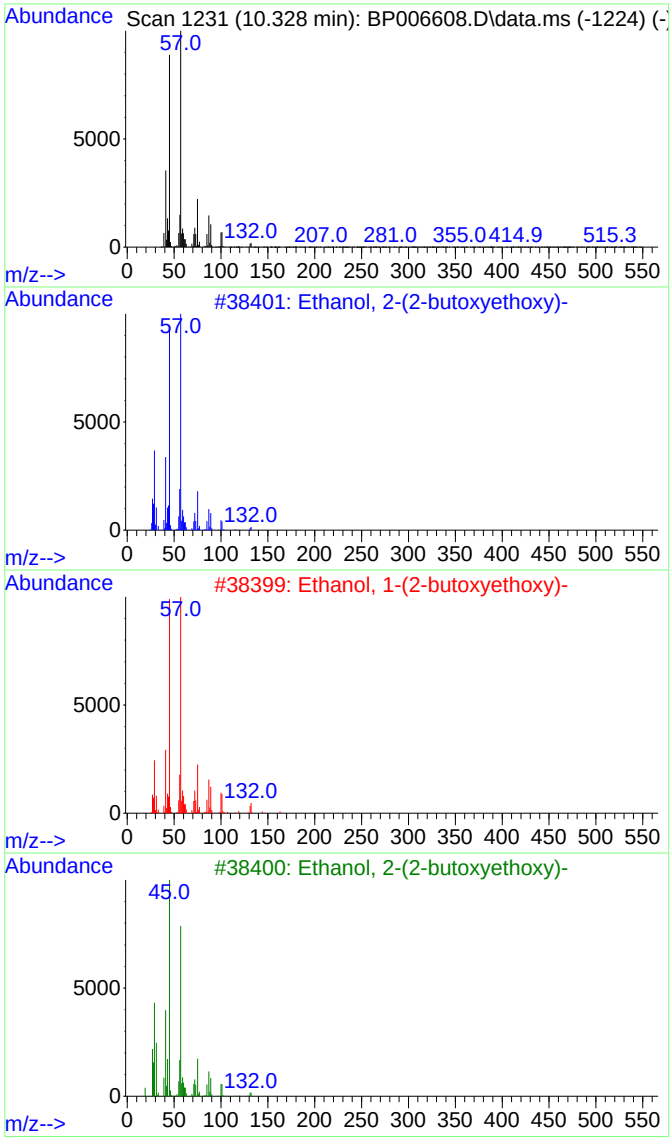
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

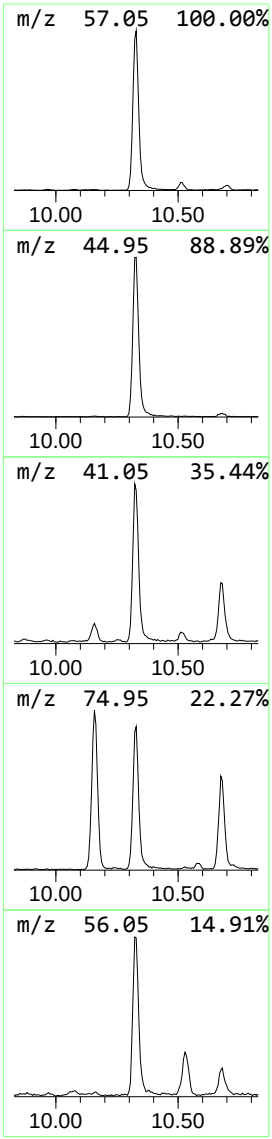
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.330	11.82 ng/ul	844300	Naphthalene-d8	10.534

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





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

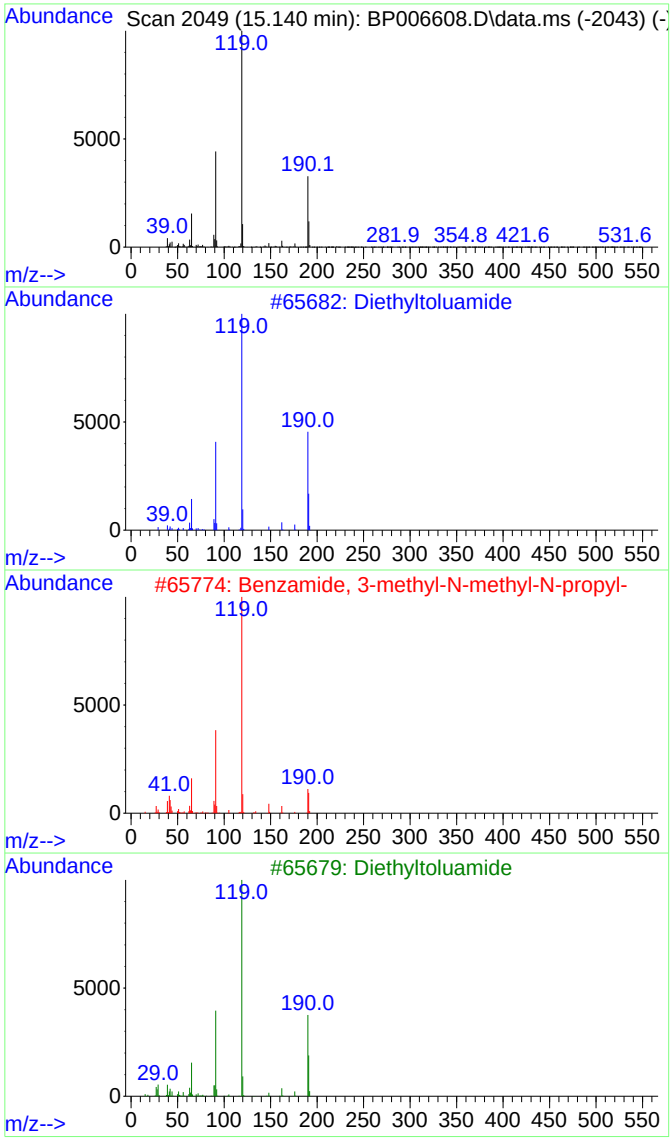
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

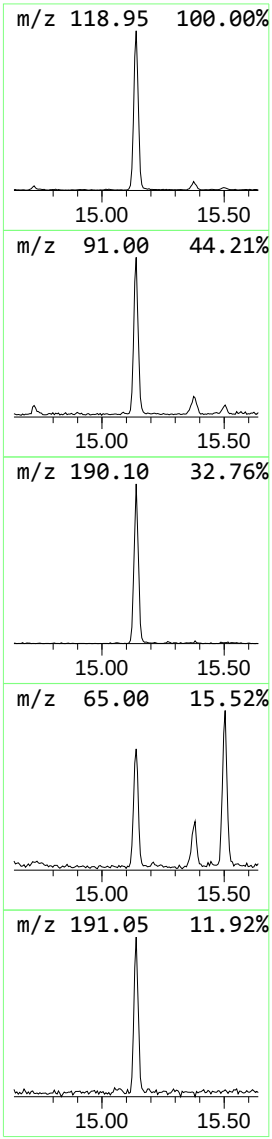
****

Peak Number 3 Diethyltoluamide Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.140	2.07 ng/ul	178309	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	95
2	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	90
3	Diethyltoluamide	191	C12H17NO	000134-62-3	90
4	Benzamide, 2-methyl-N-methyl-N-p...	191	C12H17NO	1000421-44-4	86
5	Diethyltoluamide	191	C12H17NO	000134-62-3	86





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Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00A2

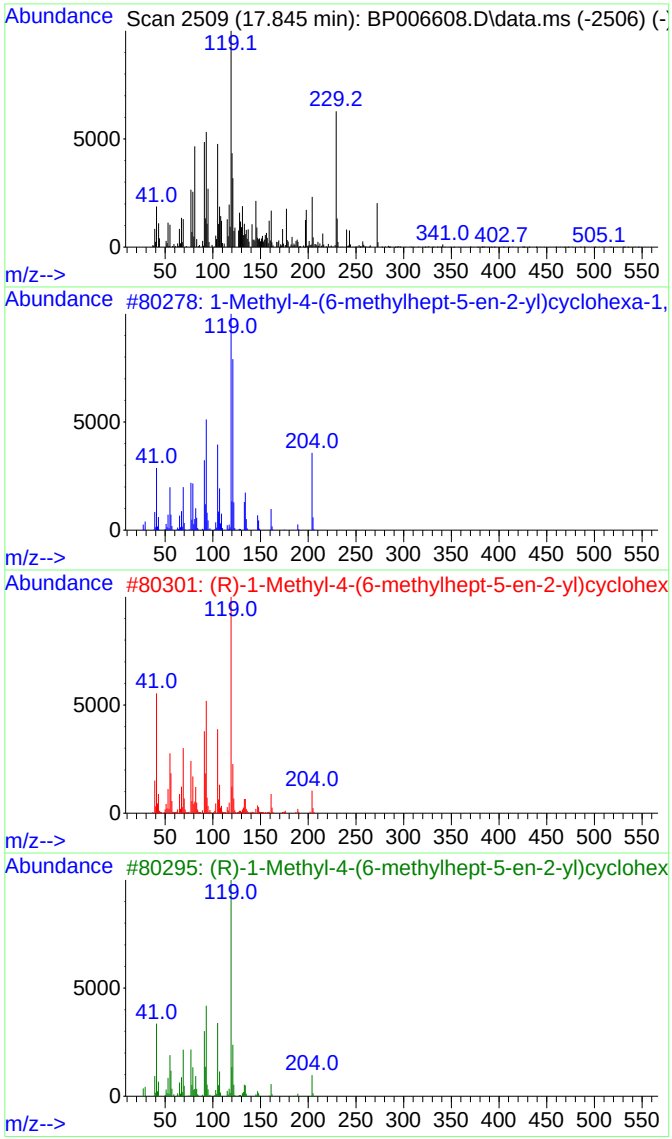
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

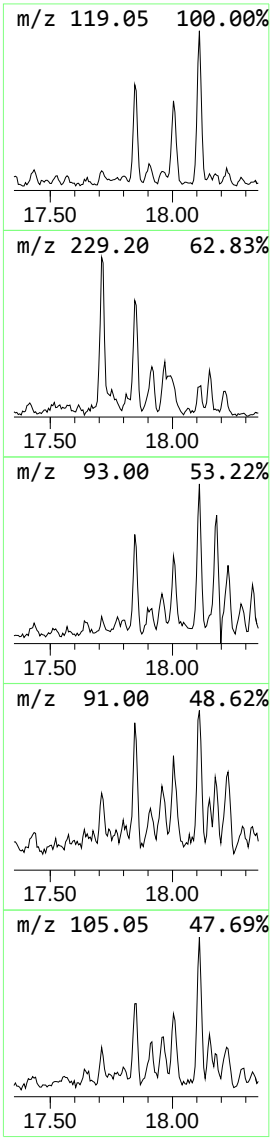
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1-Methyl-4-(6-methylhept-5-... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	2.60 ng/ul	247469	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-4-(6-methylhept-5-en-2-...	204	C15H24	000451-55-8	50
2	(R)-1-Methyl-4-(6-methylhept-5-e...	204	C15H24	028976-67-2	49
3	(R)-1-Methyl-4-(6-methylhept-5-e...	204	C15H24	028976-67-2	42
4	Di-epi-.alpha.-cedrene-(I)	204	C15H24	021996-77-0	42
5	Cubenene	204	C15H24	029837-12-5	38





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Data File : BP006608.D
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Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

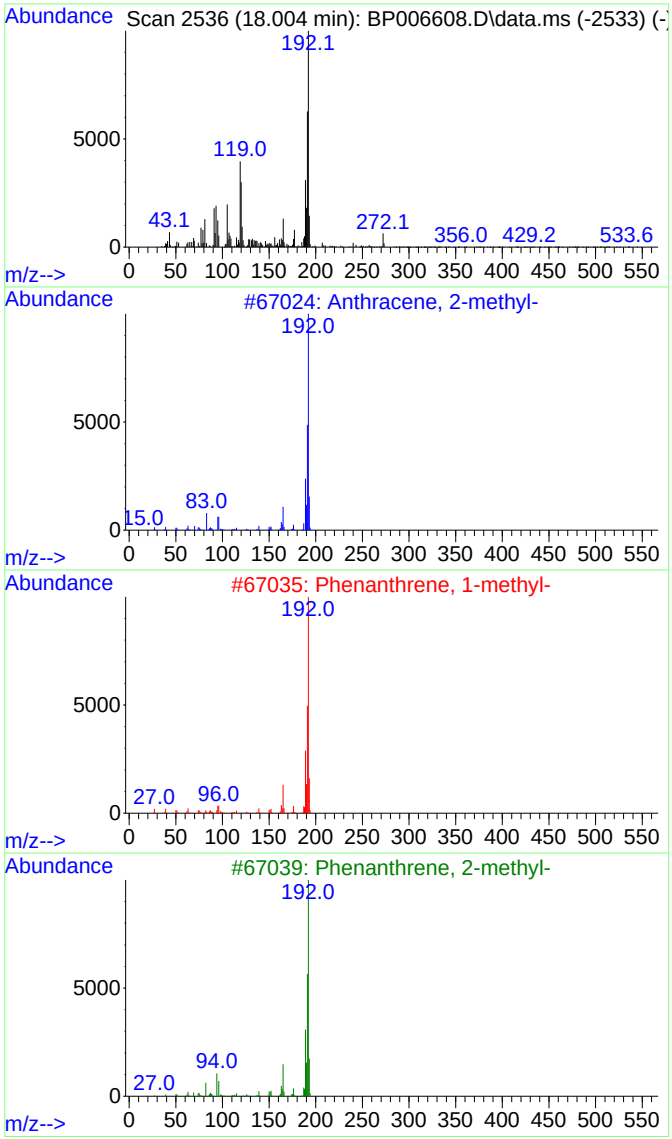
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

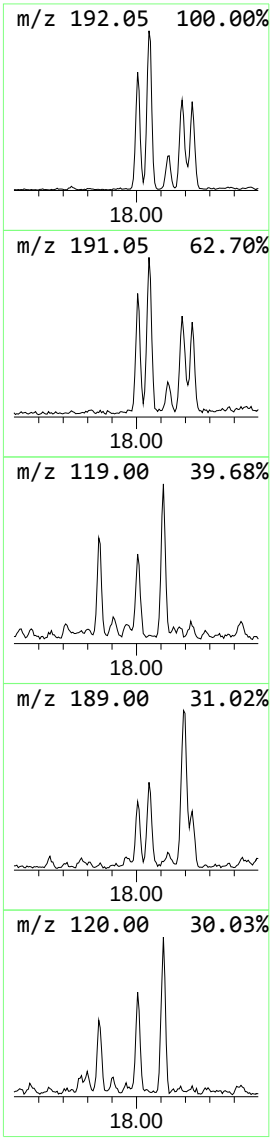
****

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.000	2.34 ng/ul	223466	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91





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Operator : CG/JU
Sample : M3169-01
Misc :
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Instrument : BNA_P
ClientSampleId : C00A2

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Quant Title : SVOA CALIBRATION

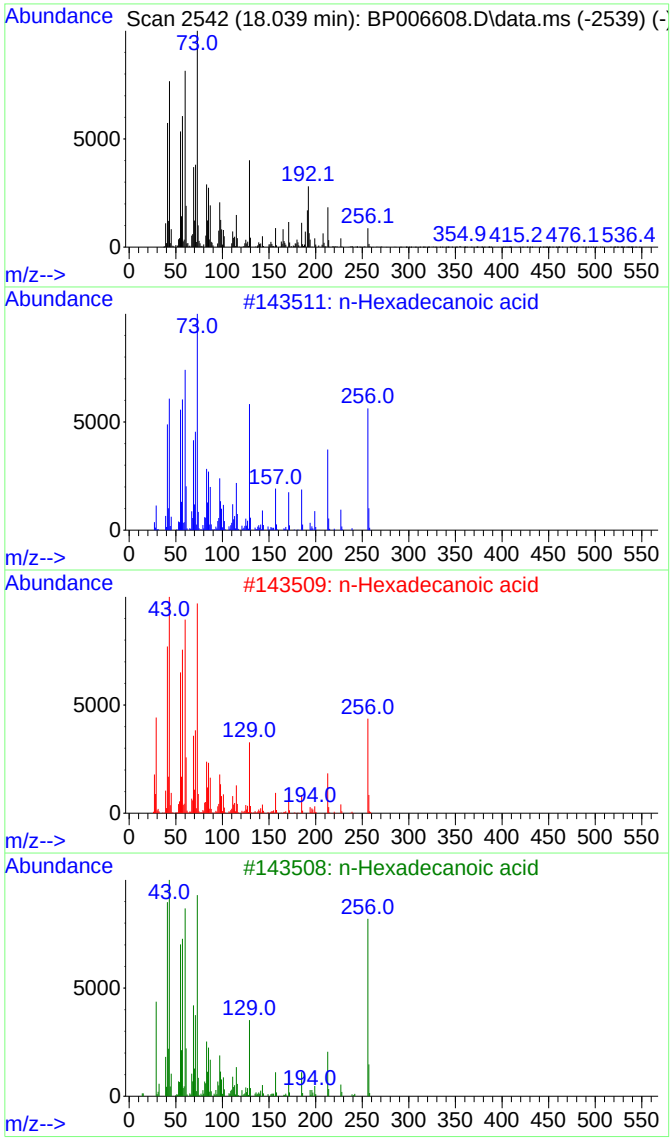
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

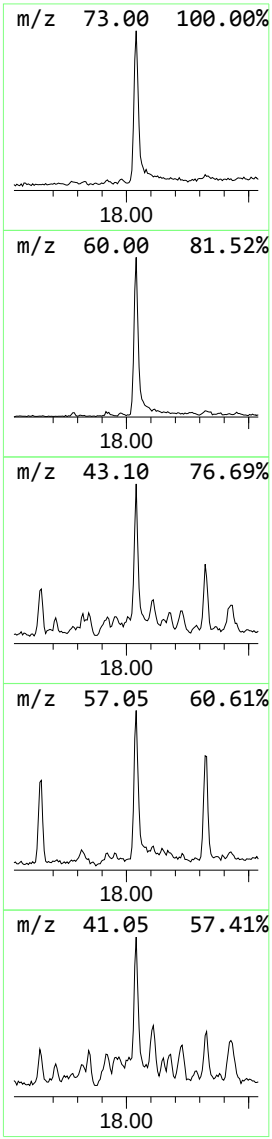
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.040	5.32 ng/ul	507068	Phenanthrene-d10	17.134

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	97
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5	Tridecanoic acid	214	C13H26O2	000638-53-9	91





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

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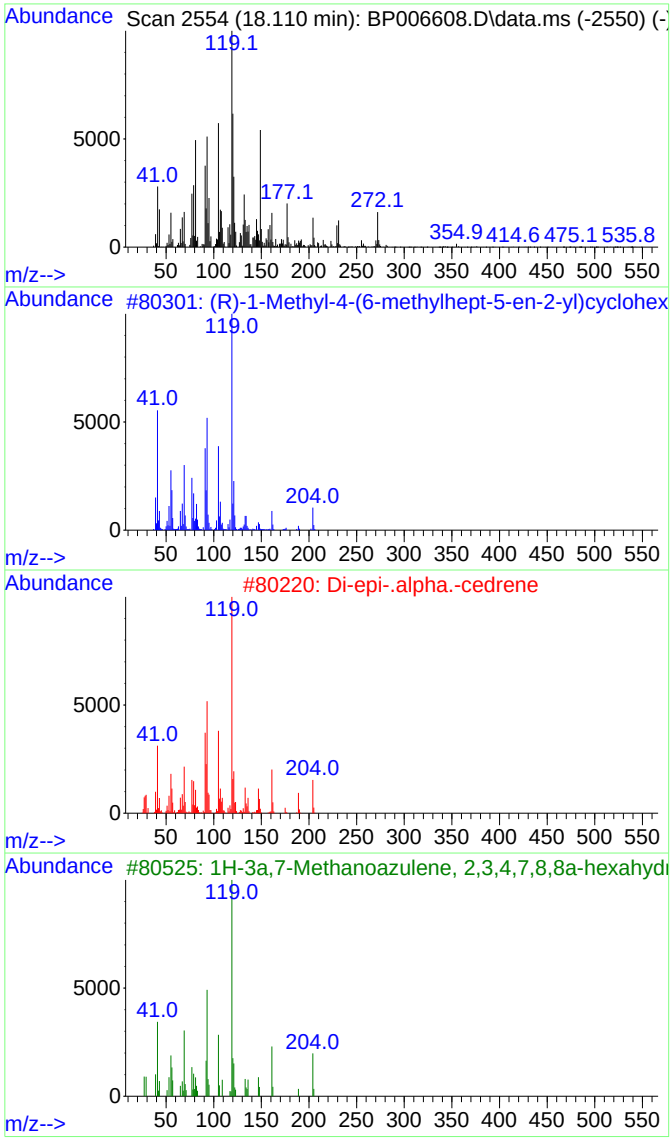
Quant Title : SVOA CALIBRATION

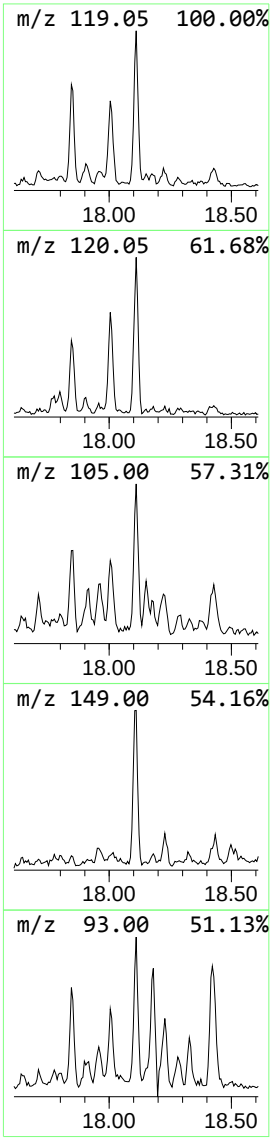
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (R)-1-Methyl-4-(6-methylhep... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.110	3.23 ng/ul	307857	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(R)-1-Methyl-4-(6-methylhept-5-e...	204	C15H24	028976-67-2	94
2	Di-epi-.alpha.-cedrene	204	C15H24	050894-66-1	62
3	1H-3a,7-Methanoazulene, 2,3,4,7,...	204	C15H24	000469-61-4	46
4	Di-epi-.alpha.-cedrene-(I)	204	C15H24	021996-77-0	46
5	(R)-1-Methyl-4-(6-methylhept-5-e...	204	C15H24	028976-67-2	46





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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

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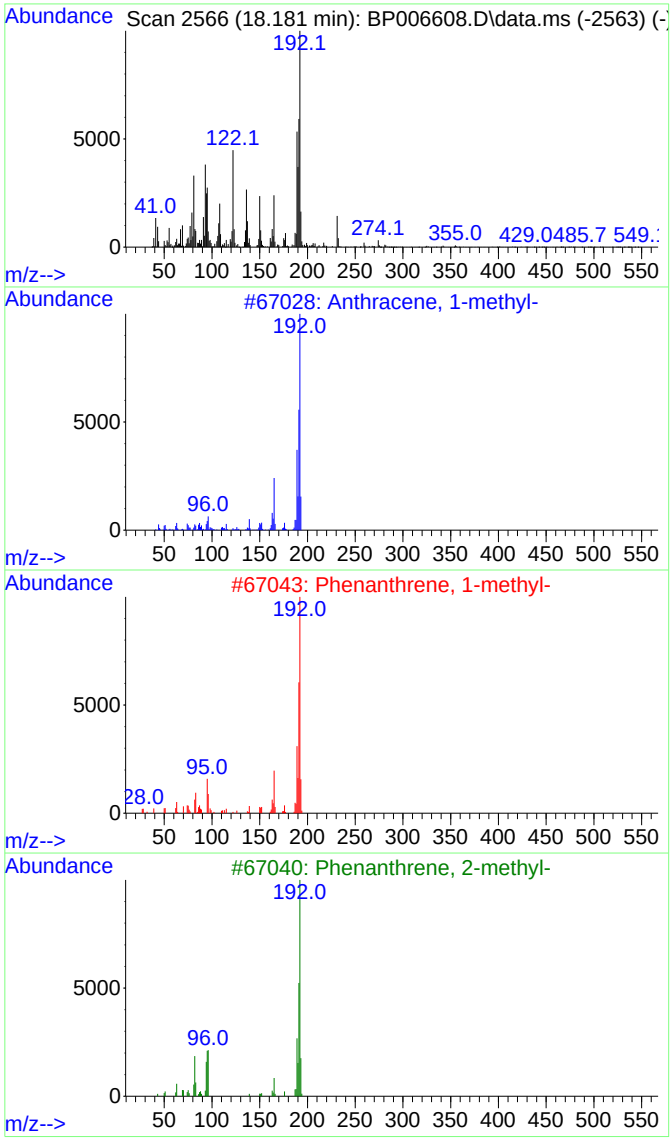
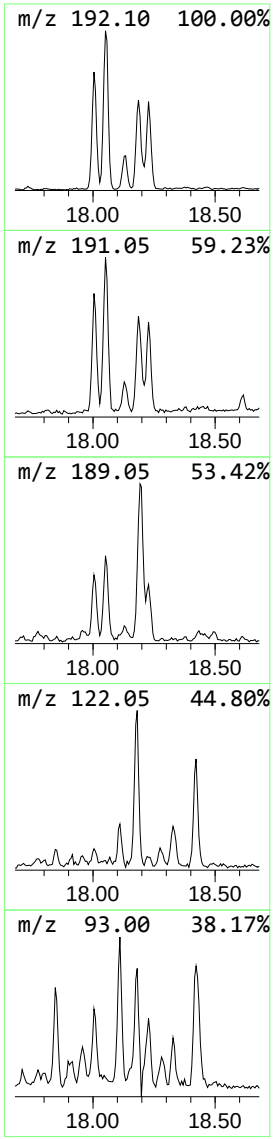
Peak Number 8 Anthracene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	2.68 ng/ul	255657	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	86
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	55
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	50





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Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

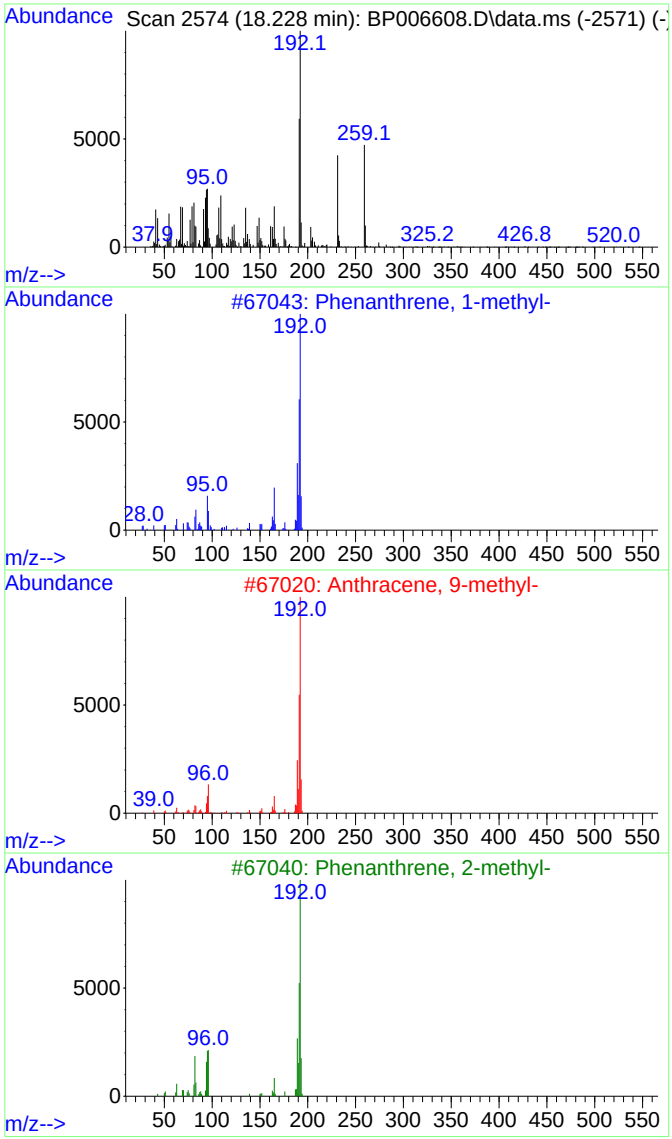
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

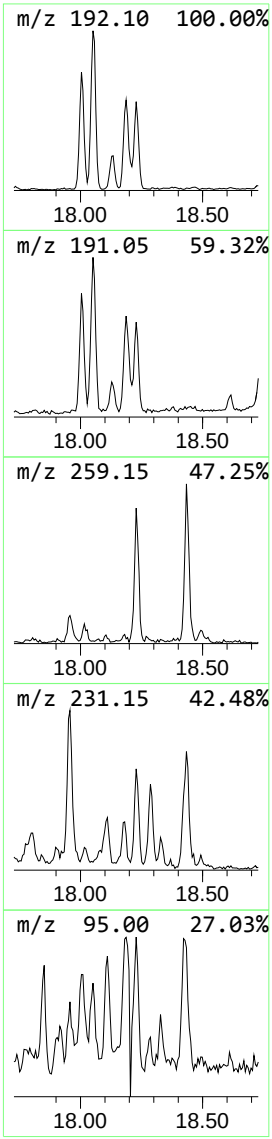
****

Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.230	2.62 ng/ul	249744	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	46
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	45
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	43
5	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	43





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Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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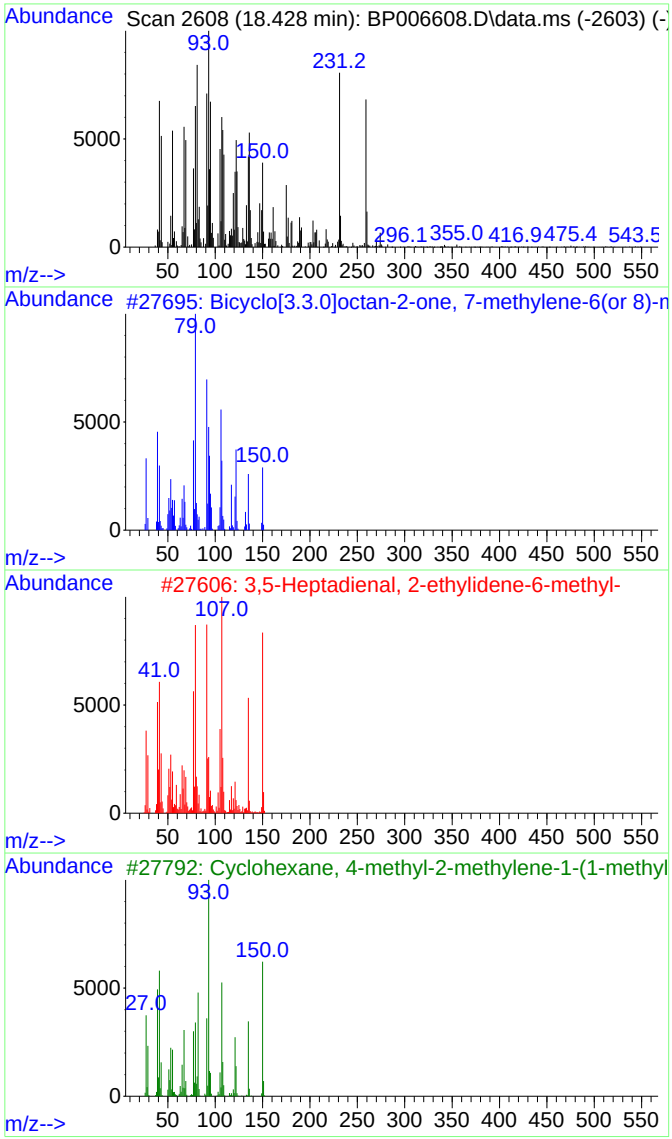
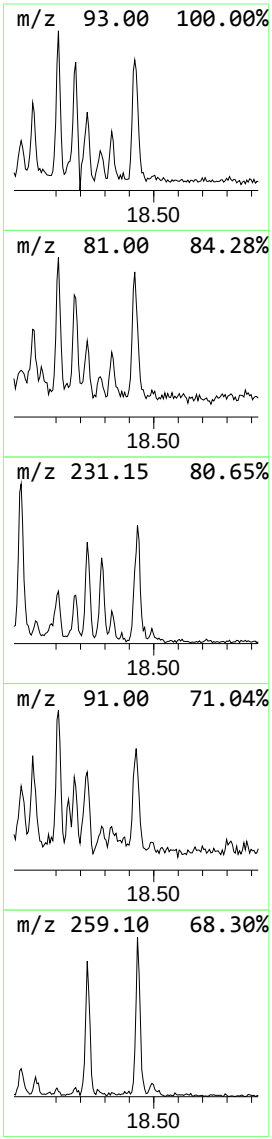
Peak Number 10 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.430	3.21 ng/ul	305878	Phenanthrene-d10	17.134

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.3.0]octan-2-one, 7-met...	150	C10H14O	1000154-23-1	25
2	3,5-Heptadienal, 2-ethylidene-6-...	150	C10H14O	099172-18-6	25
3	Cyclohexane, 4-methyl-2-methylen...	150	C11H18	056763-62-3	22
4	2H-Inden-2-one, 1,4,5,6,7,7a-hex...	136	C9H12O	039163-29-6	22
5	Cyclohexene, 2,4-dimethyl-1-(1-m...	150	C11H18	056763-60-1	22





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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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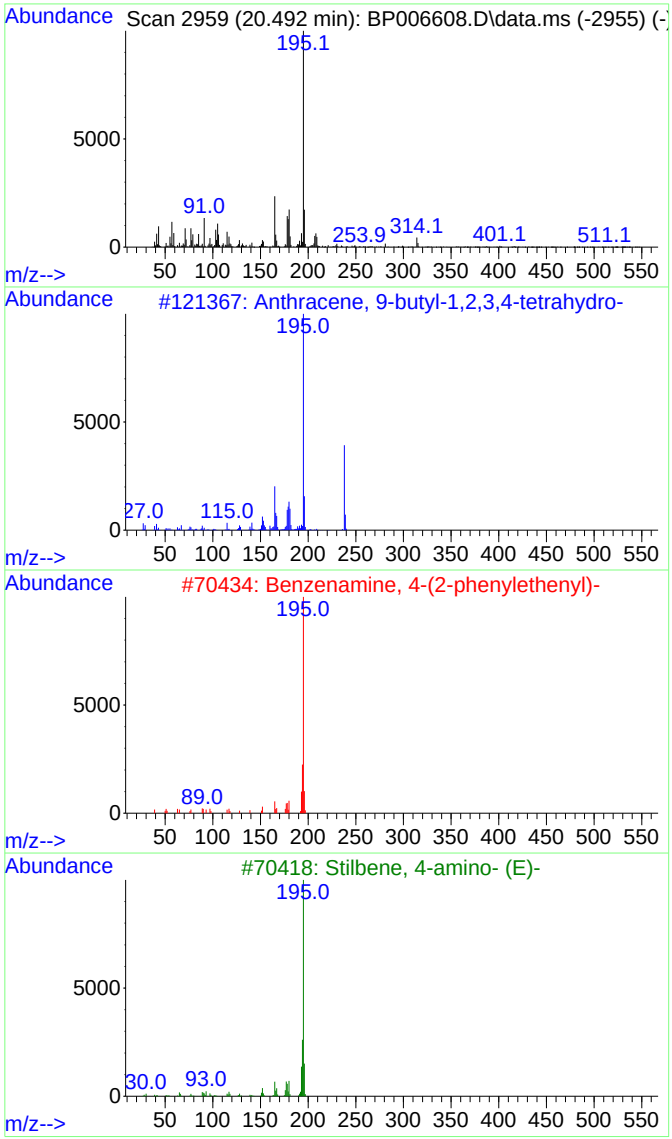
Quant Title : SVOA CALIBRATION

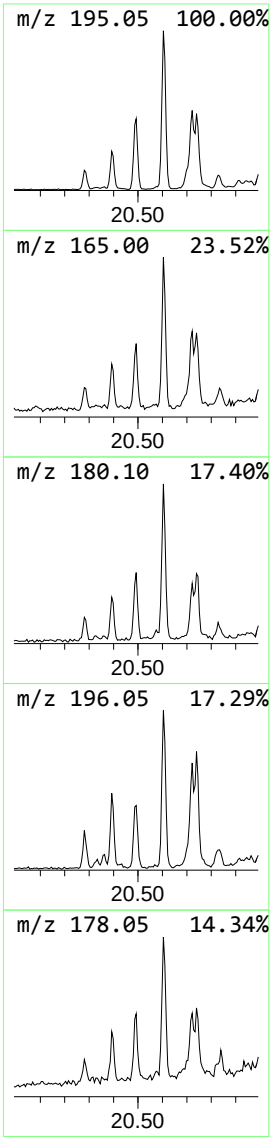
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Anthracene, 9-butyl-1,2,3,4... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.490	2.25 ng/ul	372619	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-butyl-1,2,3,4-tetr...	238	C18H22	1010151-39-2	78
2	Benzenamine, 4-(2-phenylethenyl)-	195	C14H13N	000834-24-2	72
3	Stilbene, 4-amino- (E)-	195	C14H13N	004309-66-4	72
4	9-Anthracenamine, 9,10-dihydro-	195	C14H13N	097825-91-7	64
5	Carbazole, 3,6-dimethyl-	195	C14H13N	005599-50-8	59





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Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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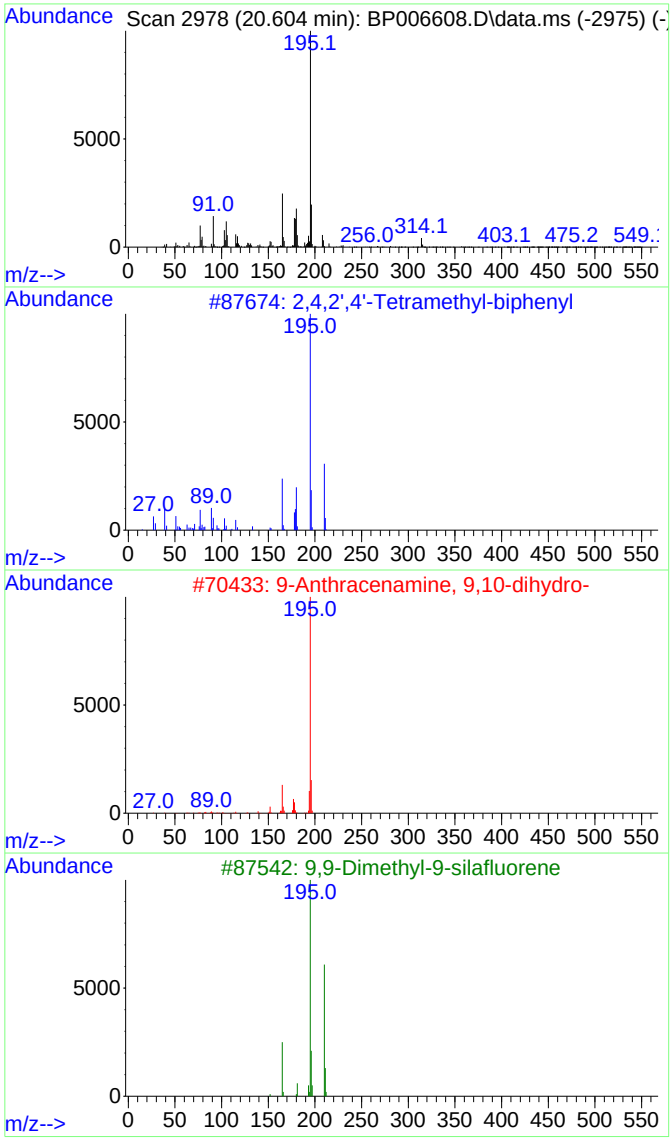
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

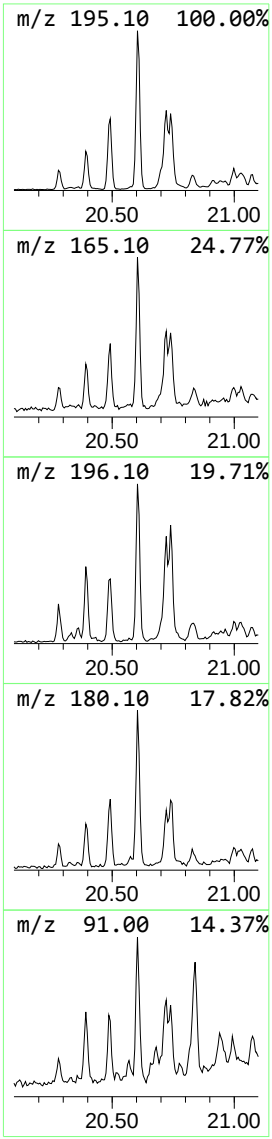
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2,4,2',4'-Tetramethyl-biphenyl Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.600	2.38 ng/ul	395131	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,2',4'-Tetramethyl-biphenyl	210	C16H18	1000486-97-7	64
2	9-Anthracenamine, 9,10-dihydro-	195	C14H13N	097825-91-7	59
3	9,9-Dimethyl-9-silafluorene	210	C14H14Si	013688-68-1	59
4	Stilbene, 4-amino- (E)-	195	C14H13N	004309-66-4	58
5	1H-Benz[f]isoindol-3-amine, 1-im...	195	C12H9N3	065558-69-2	53





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Instrument : BNA_P
ClientSampleId : C00A2

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Quant Title : SVOA CALIBRATION

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

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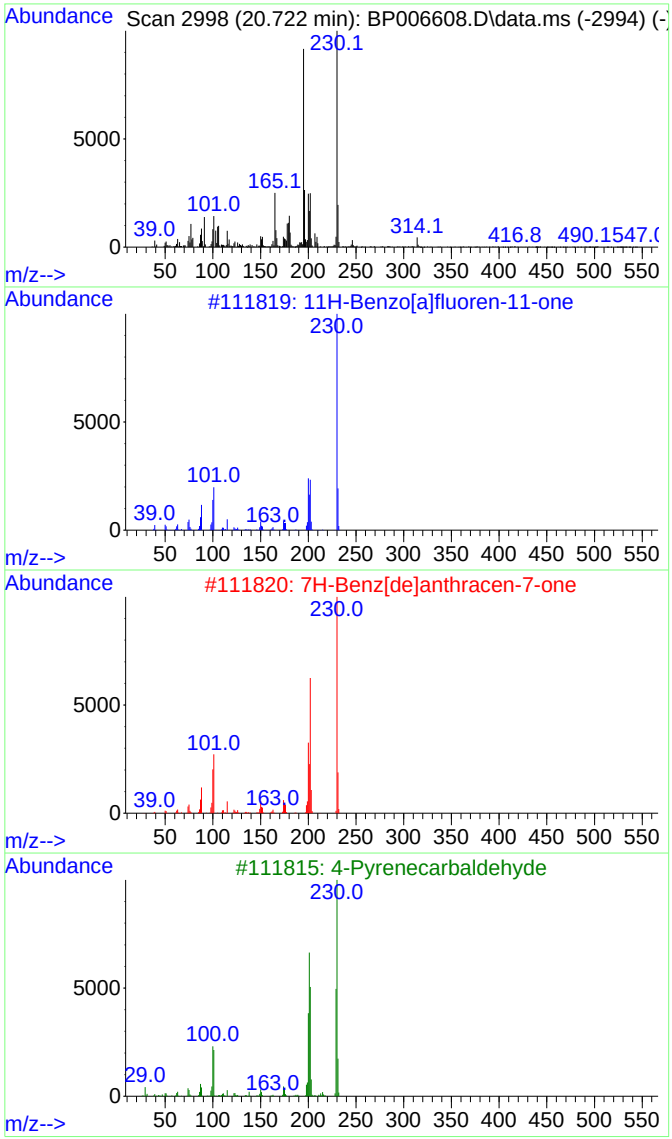
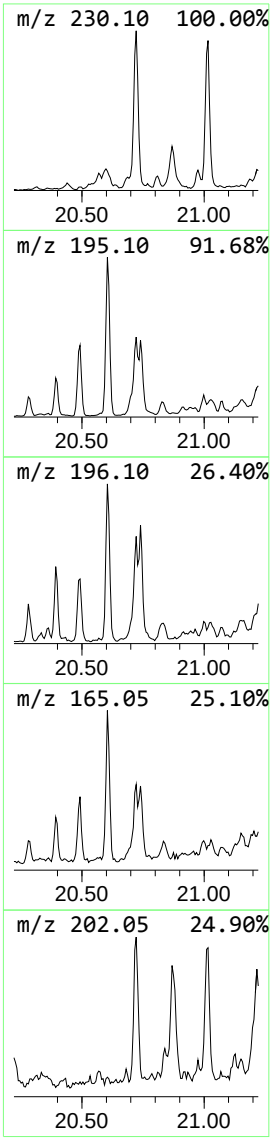
Peak Number 13 11H-Benzo[a]fluoren-11-one Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.720	3.15 ng/ul	522207	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	95
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	50
3	4-Pyrenecarbaldehyde	230	C17H10O	022245-51-8	45
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	41
5	6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	41





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Instrument : BNA_P
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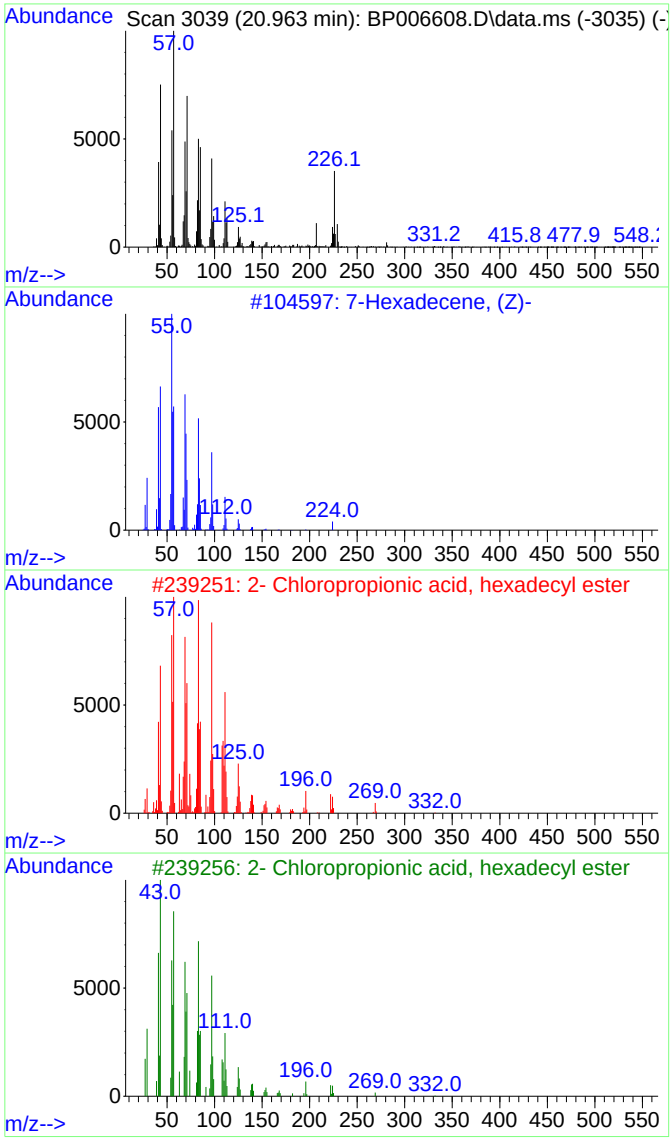
Quant Title : SVOA CALIBRATION

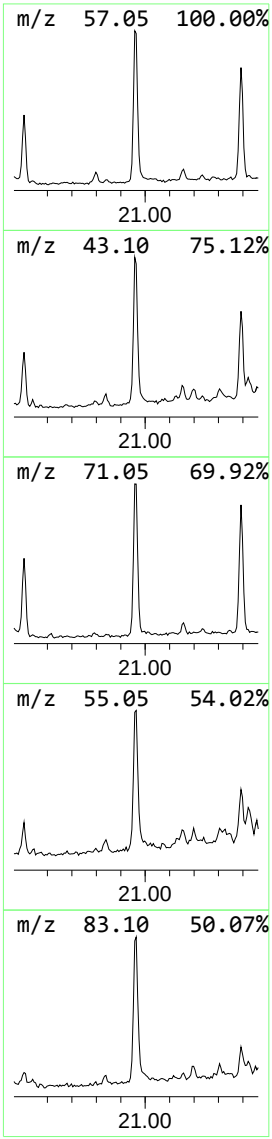
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.960	4.79 ng/ul	794700	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Hexadecene, (Z)-	224	C16H32	035507-09-6	90
2	2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	78
3	2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	78
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	70
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	70





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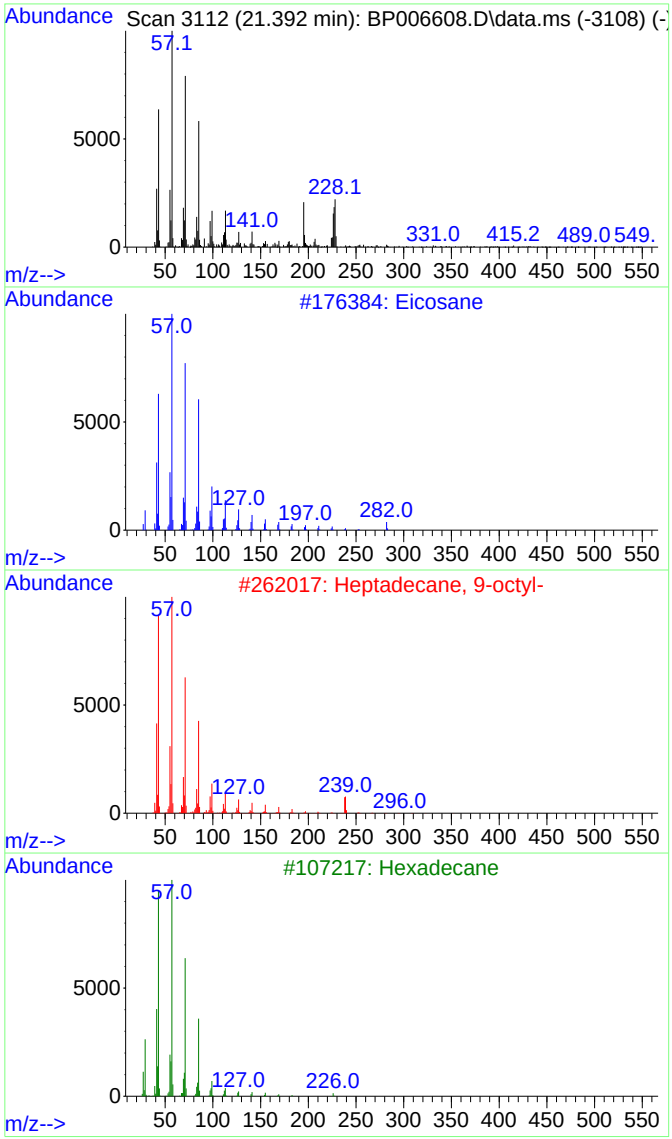
Quant Title : SVOA CALIBRATION

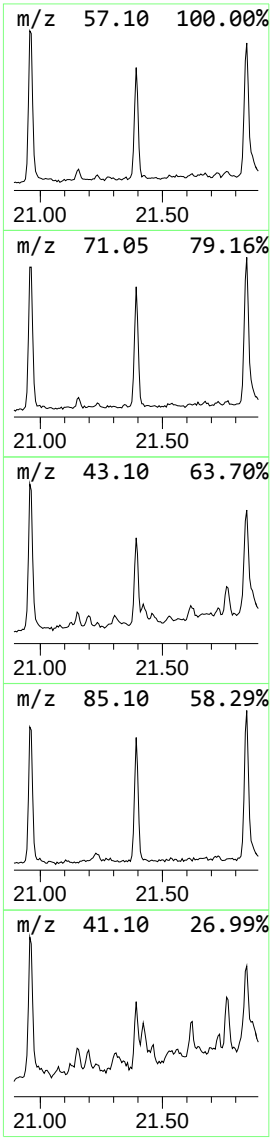
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.390	2.75 ng/ul	455963	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	78
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	78
3	Hexadecane	226	C16H34	000544-76-3	64
4	Heneicosane	296	C21H44	000629-94-7	60
5	Tetracosane	338	C24H50	000646-31-1	58





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Quant Title : SVOA CALIBRATION

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

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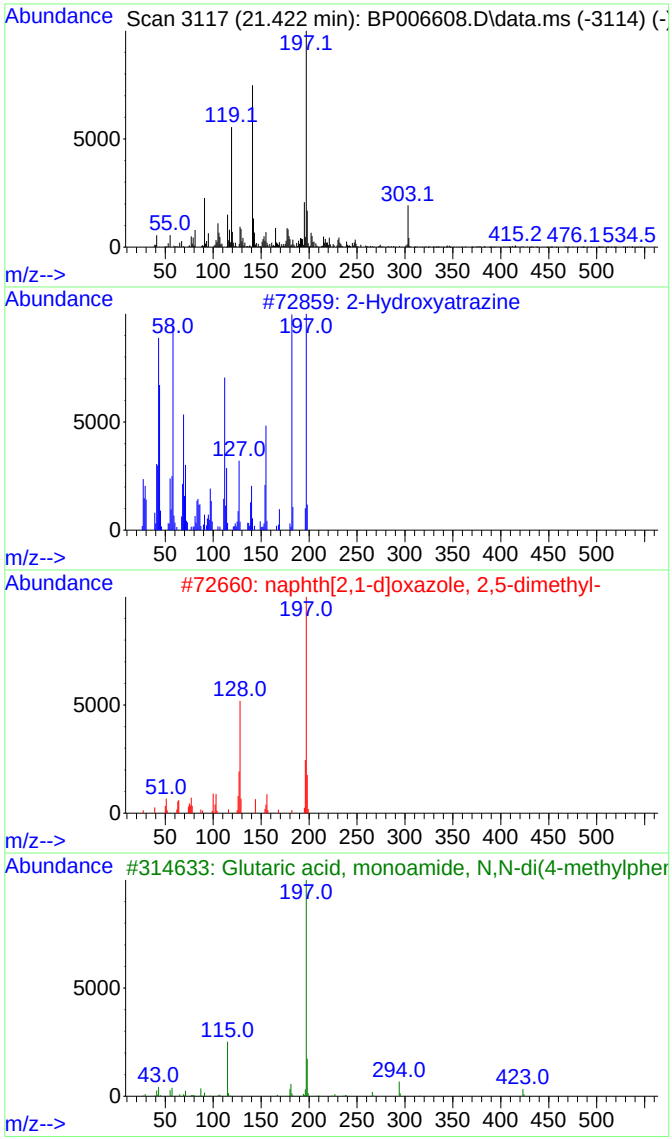
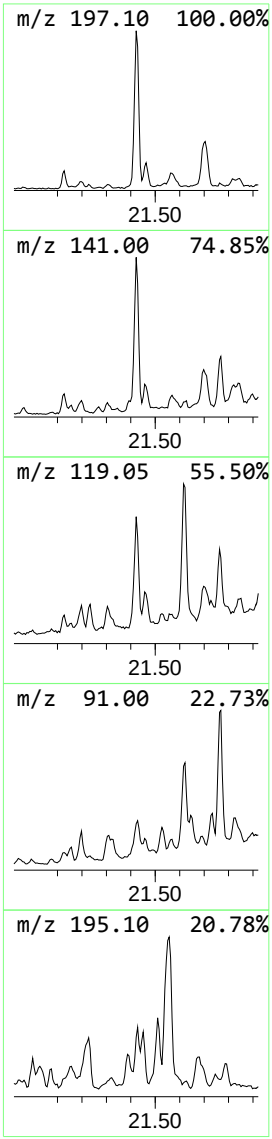
Peak Number 16 unknown-02 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.420	2.77 ng/ul	460127	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxyatrazine	197	C8H15N5O	002163-68-0	30
2	naphth[2,1-d]oxazole, 2,5-dimethyl-	197	C13H11NO	1010396-01-9	25
3	Glutaric acid, monoamide, N,N-di...	423	C27H37NO3	1000360-23-7	22
4	Glutaric acid, monoamide, N,N-di...	395	C25H33NO3	1000360-23-4	22
5	1-(4-Phenoxyphenyl)-2-phenyletha...	288	C20H16O2	003669-48-5	22





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Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00A2

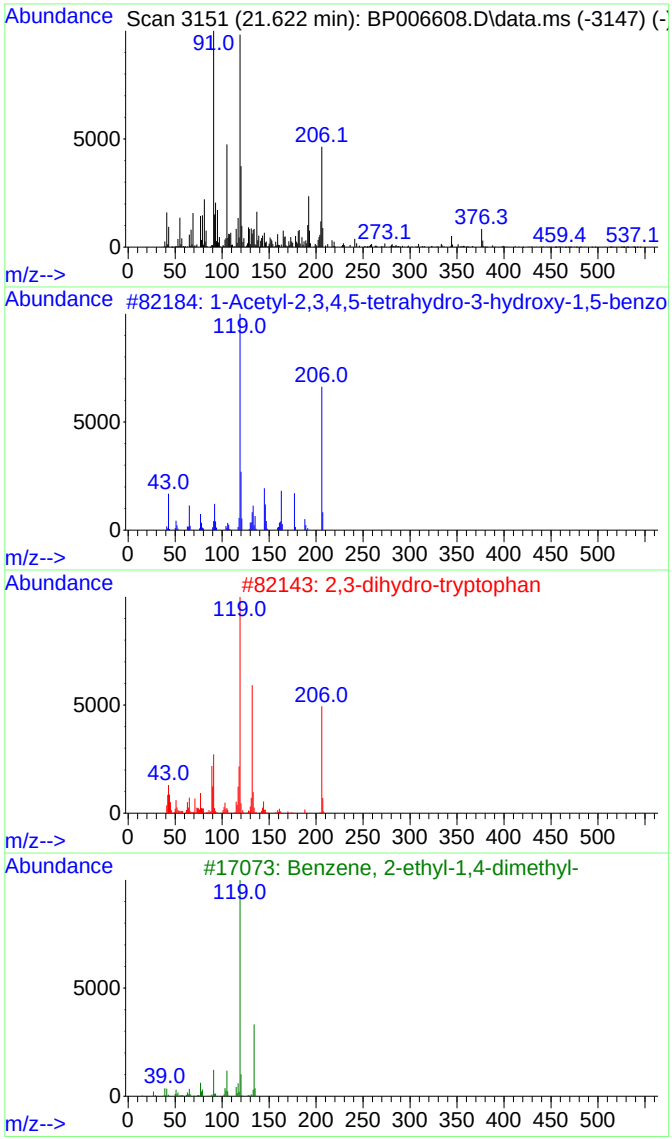
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

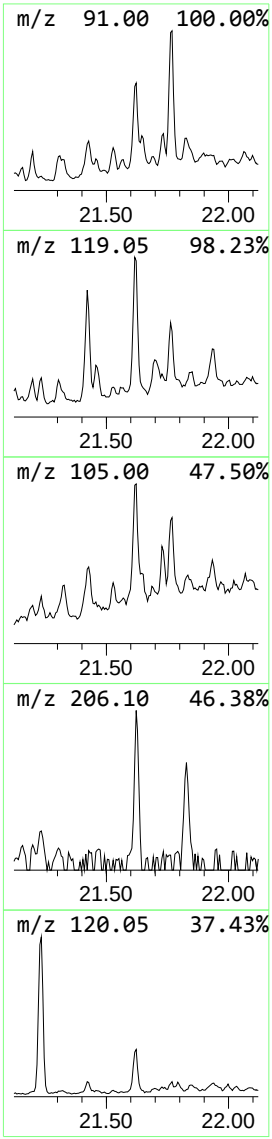
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 1-Acetyl-2,3,4,5-tetrahydro... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.620	3.13 ng/ul	520101	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Acetyl-2,3,4,5-tetrahydro-3-hy...	206	C11H14N2O2	138792-63-9	53
2	2,3-dihydro-tryptophan	206	C11H14N2O2	1000130-45-8	49
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	38
4	p-Toluic acid, undec-2-enyl ester	288	C19H28O2	1000292-50-8	35
5	1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	35





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Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00A2

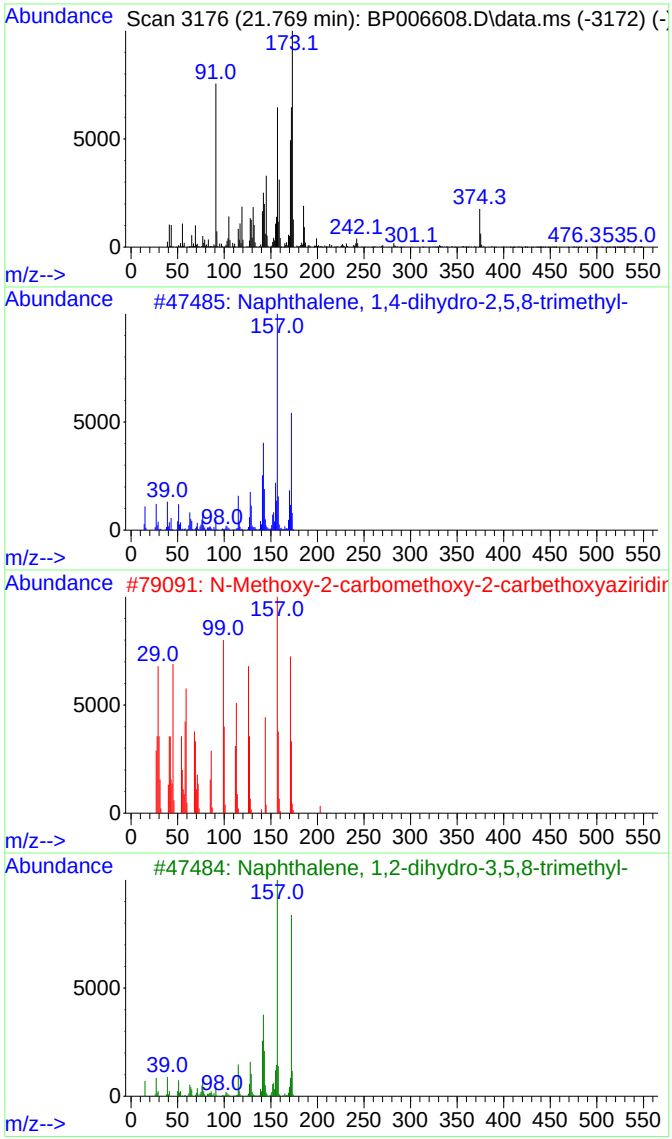
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

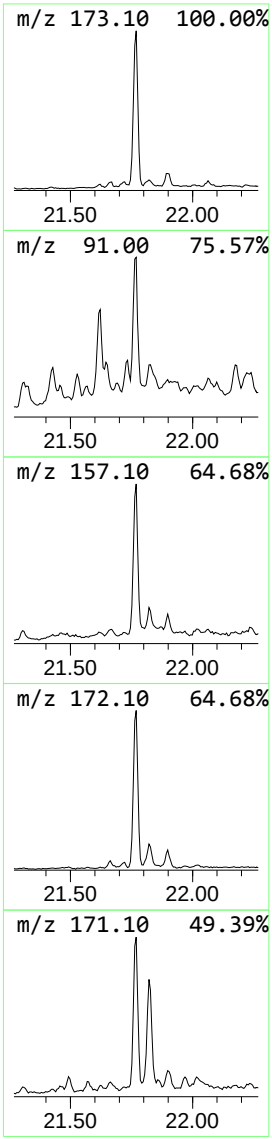
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Naphthalene, 1,4-dihydro-2,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.770	6.25 ng/ul	1036960	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4-dihydro-2,5,8-t...	172	C13H16	030316-19-9	78
2	N-Methoxy-2-carbomethoxy-2-carbe...	203	C8H13NO5	063859-01-8	70
3	Naphthalene, 1,2-dihydro-3,5,8-t...	172	C13H16	030316-18-8	70
4	1,2,3,5,6,7-Hexahydro-s-indacen-...	173	C12H15N	063089-56-5	70
5	2-(Trifluoromethyl)benzoic acid....	316	C17H23F3O2	1000283-02-0	43





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Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

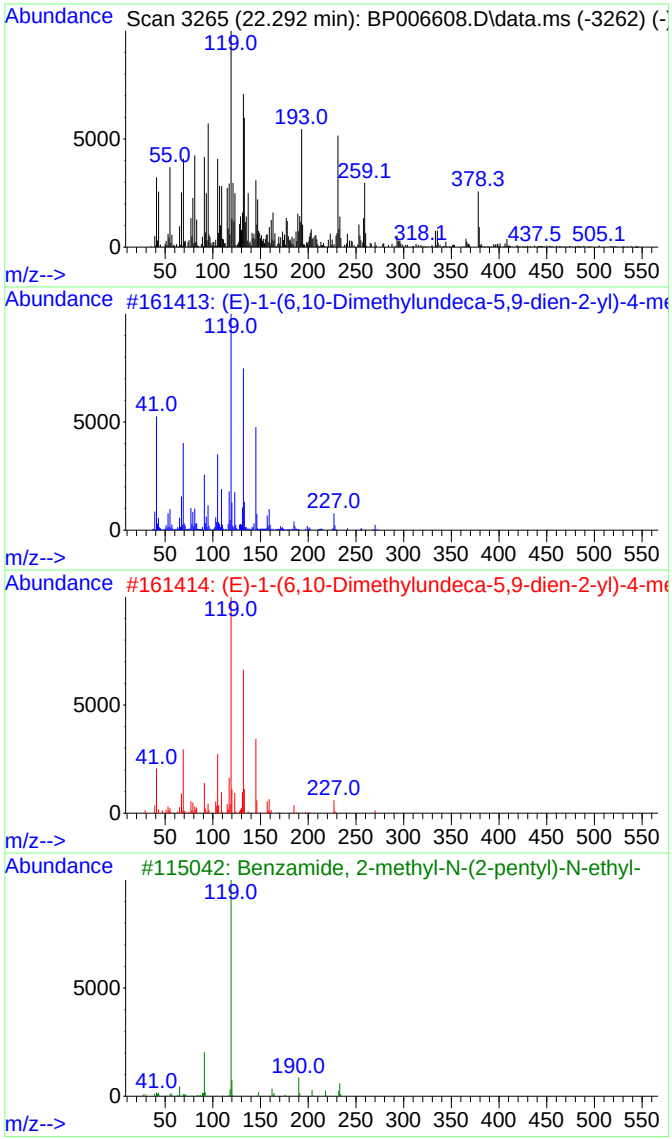
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

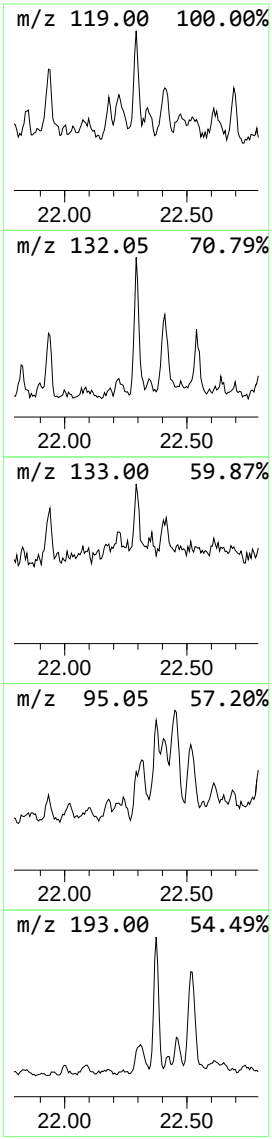
****

Peak Number 19 unknown-03 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.290	2.54 ng/ul	420756	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-(6,10-Dimethylundeca-5,9-d...	270	C20H30	055968-43-9	42
2	(E)-1-(6,10-Dimethylundeca-5,9-d...	270	C20H30	055968-43-9	38
3	Benzamide, 2-methyl-N-(2-pentyl)...	233	C15H23NO	1000490-61-0	25
4	Pentamethylnitrobenzene	193	C11H15NO2	013171-59-0	25
5	Benzamide, 4-methyl-N-butyl-N-pr...	233	C15H23NO	1000415-91-9	25





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Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00A2

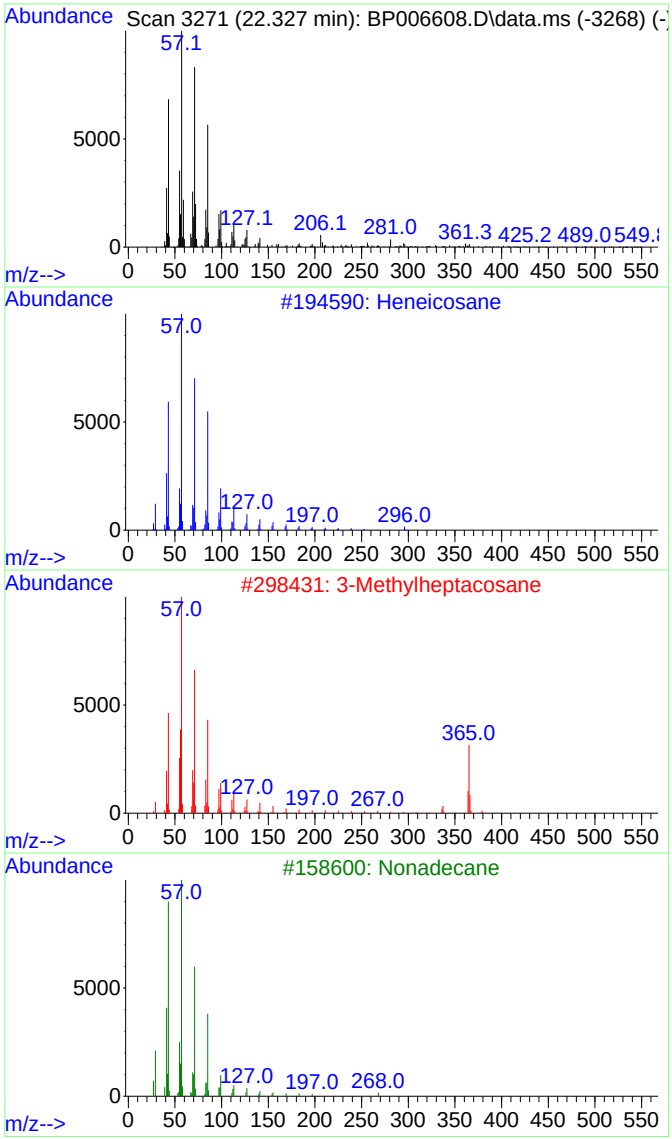
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

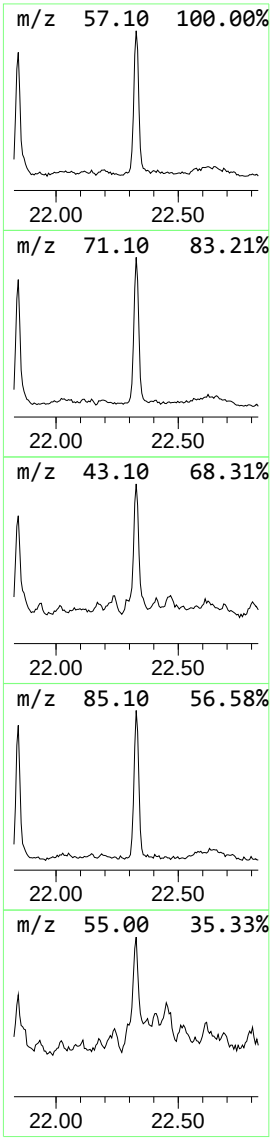
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.330	4.90 ng/ul	813372	Chrysene-d12	21.233

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	94
2	3-Methylheptacosane	394	C28H58	014167-66-9	89
3	Nonadecane	268	C19H40	000629-92-5	89
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5	Tetracosane	338	C24H50	000646-31-1	76





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Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

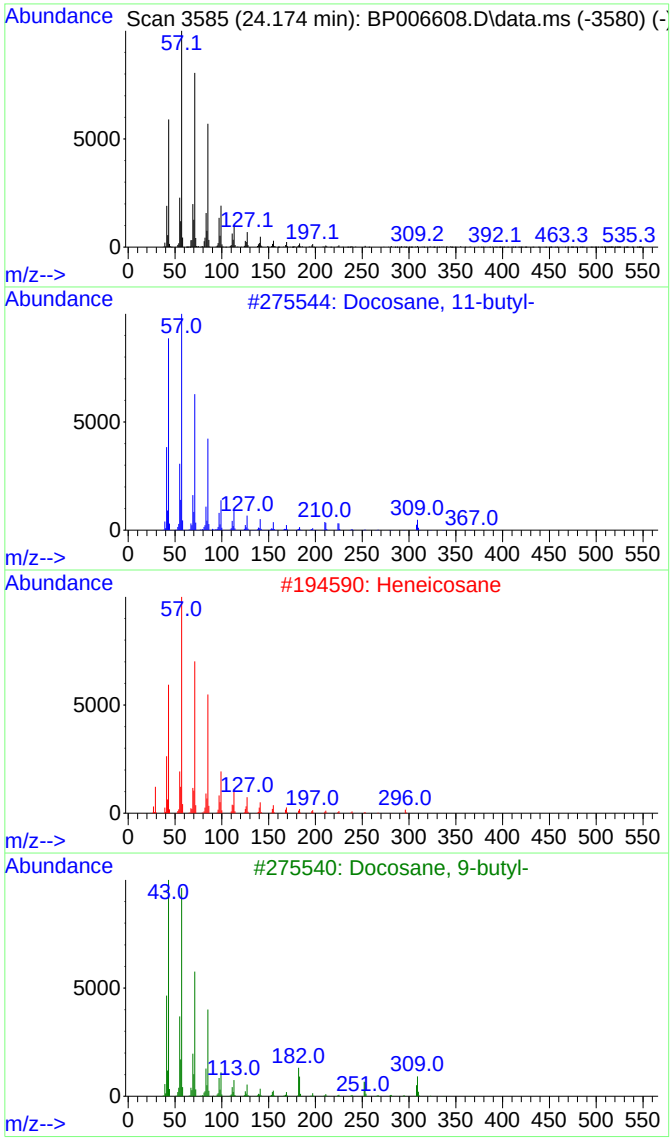
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

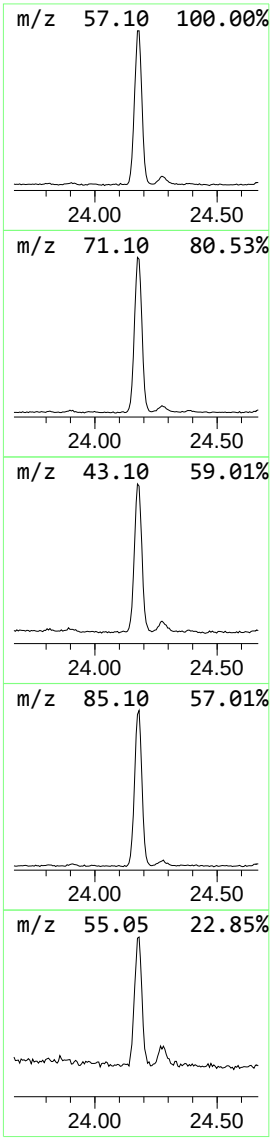
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.170	19.01 ng/ul	1792310	Perylene-d12	23.563

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Docosane, 9-butyl-	366	C26H54	055282-14-9	91
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
5	Docosane, 5-butyl-	366	C26H54	055282-16-1	87





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Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C00A2

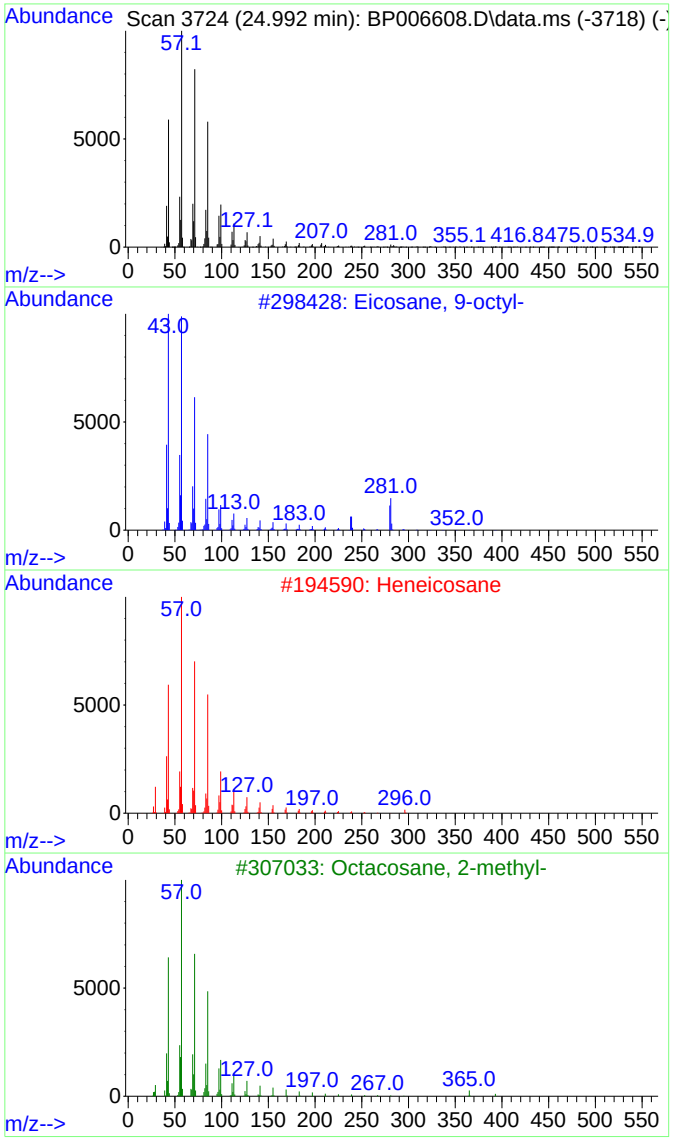
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

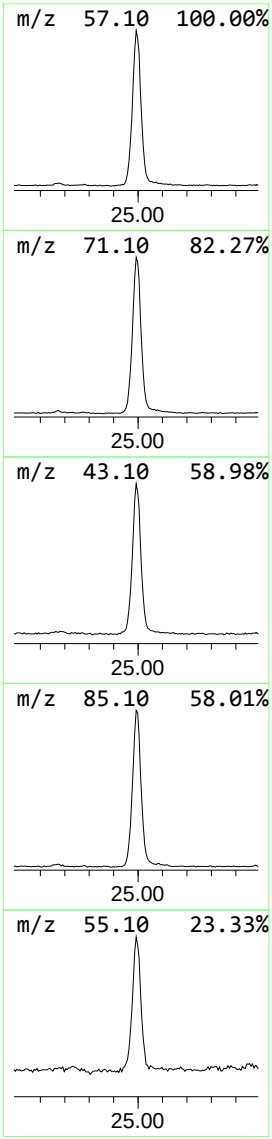
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.990	23.05 ng/ul	2172710	Perylene-d12	23.563

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
2	Heneicosane	296	C21H44	000629-94-7	91
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	90
5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

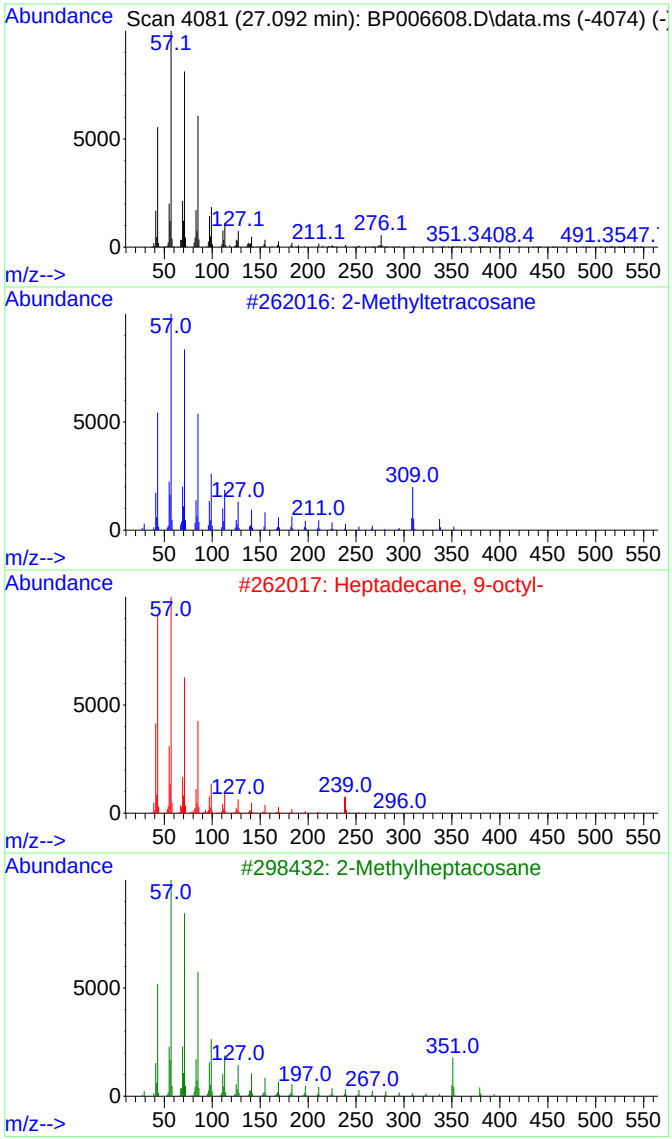
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

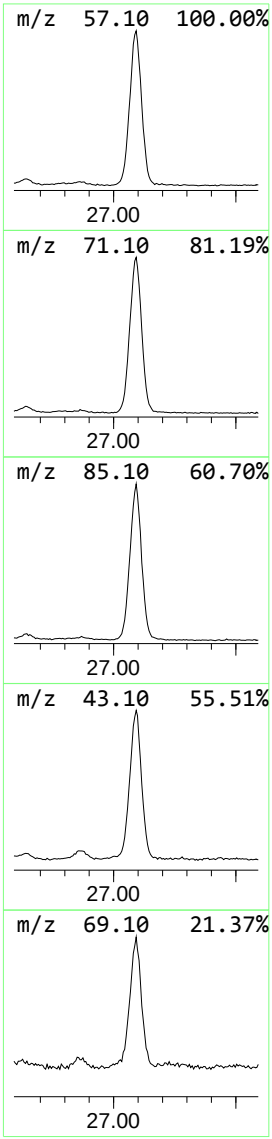
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.090	33.81 ng/ul	3186730	Perylene-d12	23.563

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyltetracosane	352	C25H52	001560-78-7	95
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3	2-Methylheptacosane	394	C28H58	001561-00-8	94
4	Heneicosane	296	C21H44	000629-94-7	91
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91





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Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

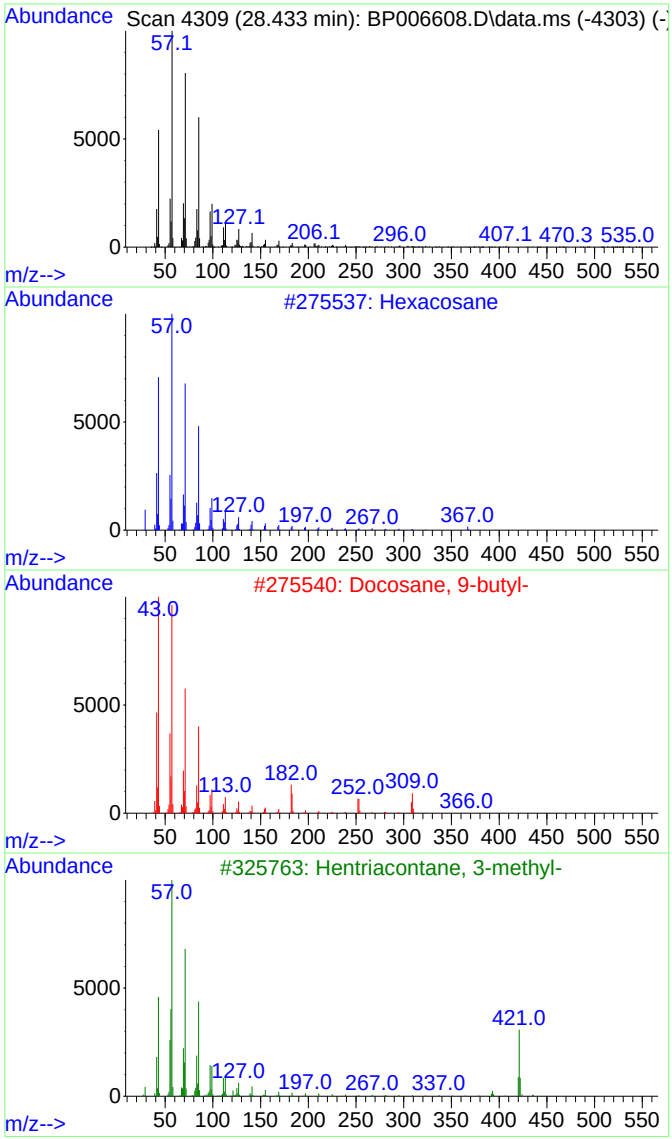
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

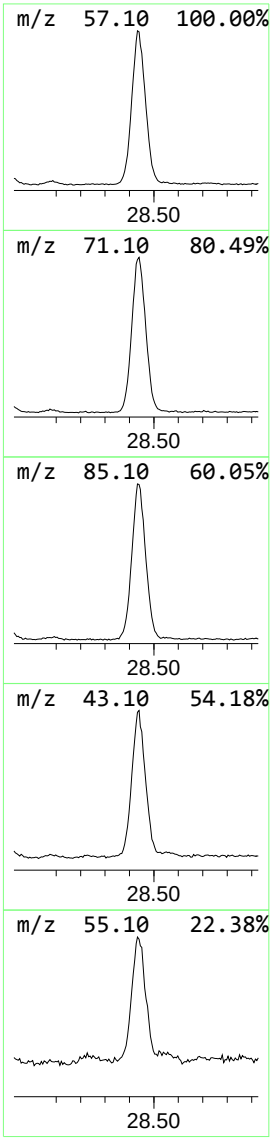
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.430	27.05 ng/ul	2550230	Perylene-d12	23.563

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Docosane, 9-butyl-	366	C26H54	055282-14-9	90
3	Hentriacontane, 3-methyl-	451	C32H66	004981-99-1	90
4	2-Methyltetracosane	352	C25H52	001560-78-7	90
5	Heneicosane	296	C21H44	000629-94-7	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080921\
Data File : BP006608.D
Acq On : 09 Aug 2021 11:26
Operator : CG/JU
Sample : M3169-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : C00A2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.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
2(3H)-Furanone,...	3.780	2.8	ng/ul	136132	1	7.752	957322	20.0
Ethanol, 2-(2-b...	10.330	11.8	ng/ul	844300	2	10.534	1428670	20.0
Diethyltoluamide	15.140	2.1	ng/ul	178309	3	14.381	1722430	20.0
1-Methyl-4-(6-m...	17.850	2.6	ng/ul	247469	4	17.134	1906990	20.0
Anthracene, 2-m...	18.000	2.3	ng/ul	223466	4	17.134	1906990	20.0
n-Hexadecanoic ...	18.040	5.3	ng/ul	507068	4	17.134	1906990	20.0
(R)-1-Methyl-4-...	18.110	3.2	ng/ul	307857	4	17.134	1906990	20.0
Anthracene, 1-m...	18.180	2.7	ng/ul	255657	4	17.134	1906990	20.0
Phenanthrene, 1...	18.230	2.6	ng/ul	249744	4	17.134	1906990	20.0
unknown-01	18.430	3.2	ng/ul	305878	4	17.134	1906990	20.0
Anthracene, 9-b...	20.490	2.3	ng/ul	372619	5	21.233	3318170	20.0
2,4,2',4'-Tetra...	20.600	2.4	ng/ul	395131	5	21.233	3318170	20.0
11H-Benzo[a]flu...	20.720	3.1	ng/ul	522207	5	21.233	3318170	20.0
(DEL) Alkane: S...	20.960	4.8	ng/ul	794700	5	21.233	3318170	20.0
(DEL) Alkane: S...	21.390	2.8	ng/ul	455963	5	21.233	3318170	20.0
unknown-02	21.420	2.8	ng/ul	460127	5	21.233	3318170	20.0
1-Acetyl-2,3,4,...	21.620	3.1	ng/ul	520101	5	21.233	3318170	20.0
Naphthalene, 1,...	21.770	6.3	ng/ul	1036960	5	21.233	3318170	20.0
unknown-03	22.290	2.5	ng/ul	420756	5	21.233	3318170	20.0
(DEL) Alkane: S...	22.330	4.9	ng/ul	813372	5	21.233	3318170	20.0
(DEL) Alkane: S...	24.170	19.0	ng/ul	1792310	6	23.563	1885350	20.0
(DEL) Alkane: S...	24.990	23.1	ng/ul	2172710	6	23.563	1885350	20.0
(DEL) Alkane: S...	27.090	33.8	ng/ul	3186730	6	23.563	1885350	20.0
(DEL) Alkane: S...	28.430	27.1	ng/ul	2550230	6	23.563	1885350	20.0