LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDC70

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-BP092424.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP022178.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.758	465	471	480	rVB2	5348474	8399456	3.82%	0.978%
2	6.222	544	550	557	rBV5	1171378	2689751	1.22%	0.313%
3	6.287	557	561	566	rBV	2094156	2934301	1.33%	0.342%
4	6.369	566	575	578	rBV3	1717368	3858634	1.75%	0.449%
5	6.716	630	634	639	rVB2	2707941	4206703	1.91%	0.490%
6	6.769	639	643	647	rBV	3140019	4945362	2.25%	0.576%
7	6.810	647	650	655	rVB	2784900	3919932	1.78%	0.456%
8	6.875	655	661	668	rBV3	4280442	9049499	4.11%	1.053%
9	6.940	668	672	679	rVB	2493387	4024941	1.83%	0.468%
10	7.099	694	699	703	rBV2	2428480	3940129	1.79%	0.459%
11	7.146	703	707	714	rBV2	1571782	2595518	1.18%	0.302%
12	7.352	734	742	750	rVB2	14660484	28272883	12.84%	3.291%
13	7.622	778	788	793	rVV4	1188357	3663827	1.66%	0.426%
14	7.687	793	799	805	rVB2	3699954	6692747	3.04%	0.779%
15	7.787	805	816	819	rBV2	6904042	14121619	6.41%	1.644%
16	7.822	819	822	829	rVV2	2063261	3625448	1.65%	0.422%
17	7.993	848	851	858	rVB3	1698771	2930652	1.33%	0.341%
18	8.134	869	875	879	rBV	1950369	3335082	1.51%	0.388%
19	8.234	888	892	898	rVB2	3035696	5768093	2.62%	0.671%
20	8.293	898	902	905	rVB	2006225	2574625	1.17%	0.300%
21	8.340	905	910	911	rBV2	3789579	5273575	2.40%	0.614%
22	8.463	926	931	934	rBV	3408289	5035065	2.29%	0.586%
23	8.499	934	937	945	rVB	2404701	3924674	1.78%	0.457%
24	8.646	957	962	966	rBV	2081377	3280262	1.49%	0.382%
25	8.699	966	971	978	rVB	2429911	4019810	1.83%	0.468%
26	8.799	978	988	993	rBV2	3418010	7352268	3.34%	0.856%
27	8.934	999	1011	1019	rVB	11858878	24411626	11.09%	2.841%
28	9.040	1019	1029	1035	rBV6	2379167	6944178	3.15%	0.808%
29	9.122	1036	1043	1048	rBV3	1309553	3167565	1.44%	0.369%
30	9.563	1109	1118	1122	rBV5	1730076	4014879	1.82%	0.467%
31	9.616	1122	1127	1131	rBV4	1332328	2390276	1.09%	0.278%
32	9.757	1148	1151	1157	rVB	1835738	2537416	1.15%	0.295%
33	9.998	1186	1192	1198	rVV2	2311372	4276635	1.94%	0.498%
34	10.163	1214	1220	1225	rBV2	1646230	2674341	1.21%	0.311%
35	10.363	1244	1254	1258	rBV5	1145012	3104013	1.41%	0.361%
36	10.451	1259	1269	1275	rVV2	7804340	17847167	8.11%	2.077%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

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-BP092424.MA.M
Title : SVOA CALIBRATION

37	10.510	1275	1279	1286	rVV2	1402479	2845926	1.29%	0.331%
38	10.587	1286	1292	1296	rVV4	3781390	6904141	3.14%	0.804%
39	10.840	1329	1335	1345	rBV	5646669	9324642	4.24%	1.085%
40	10.969	1348	1357	1363	rVB	1592973	3213369	1.46%	0.374%
41	11.151	1380	1388	1392	rBV5	939613	2533063	1.15%	0.295%
42	11.369	1417	1425	1434	rVB5	1195137	3257832	1.48%	0.379%
43	11.493	1439	1446	1450	rBV2	2720323	5113304	2.32%	0.595%
44	11.563	1450	1458	1464	rVB	6018222	11561265	5.25%	1.346%
45	11.722	1477	1485	1492	rVB3	1524385	3000794	1.36%	0.349%
46	11.881	1506	1512	1516	rBV	5421088	8582280	3.90%	0.999%
47	11.922	1516	1519	1523	rVB2	1930164	2583909	1.17%	0.301%
48	12.122	1547	1553	1563	rVB2	6770960	11788700	5.36%	1.372%
49	12.287	1576	1581	1585	rBV5	1372587	2653165	1.21%	0.309%
50	12.340	1585	1590	1599	rVV	2113984	4657603	2.12%	0.542%
51	12.528	1616	1622	1634	rVB5	1700643	4277929	1.94%	0.498%
52	12.698	1646	1651	1661	rBV3	1205758	2372860	1.08%	0.276%
53	13.128	1718	1724	1731	rVV	6218274	9927963	4.51%	1.156%
54	13.481	1776	1784	1794	rBV4	2542990	6540657	2.97%	0.761%
55	13.628	1803	1809	1814	rBV	3013560	6454696	2.93%	0.751%
56	13.681	1814	1818	1829	rVB2	3417557	8735203	3.97%	1.017%
57	13.792	1829	1837	1841	rBV2	2566015	4684172	2.13%	0.545%
58	13.939	1858	1862	1868	rBV2	2449710	3549182	1.61%	0.413%
59	14.234	1903	1912	1915	rBV	18147948	32862637	14.93%	3.825%
60	14.404	1936	1941	1945	rVV2	3166292	5687028	2.58%	0.662%
61	14.463	1945	1951	1955	rVV	10055041	14291742	6.49%	1.663%
62	14.528	1957	1962	1970	rVB3	2889325	6216344	2.82%	0.724%
63	14.728	1993	1996	2001	rVB2	2354595	3226018	1.47%	0.375%
64	14.798	2002	2008	2012	rBV2	2031630	3250964	1.48%	0.378%
65	14.845	2012	2016	2020	rVV2	1801997	3419500	1.55%	0.398%
66	14.981	2026	2039	2043	rVV	20680986	41562950	18.88%	4.838%
67	15.028	2043	2047	2050	rVV3	2522944	3823707	1.74%	0.445%
68	15.098	2054	2059	2066	rVV	10381996	16427603	7.46%	1.912%
69	15.186	2066	2074	2079	rVB2	6246880	12772200	5.80%	1.487%
70	15.328	2095	2098	2104	rVB	1925648	2663545	1.21%	0.310%
71	15.445	2112	2118	2124	rBV2	3402896	6089734	2.77%	0.709%
72	15.598	2139	2144	2150	rVB2	2457625	4454325	2.02%	0.518%
73	15.704	2157	2162	2167	rVB4	2762749	3823762	1.74%	0.445%
74	15.975	2203	2208	2214	rVB3	2773821	4058371	1.84%	0.472%
75	16.075	2219	2225	2228	rBV2	7796709	13561115	6.16%	1.578%
76	16.816	2337	2351	2360	rBV5	38612043	220138665	100.00%	25.623%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDC70

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-BP092424.MA.M
Title : SVOA CALIBRATION

77	16.933	2367	2371	2383	rVB2	4098276	8330524	3.78%	0.970%
78	17.145	2401	2407	2411	rVV3	2874057	5801302	2.64%	0.675%
79	17.180	2411	2413	2417	rVV	1768168	2895109	1.32%	0.337%
80	17.228	2417	2421	2425	rVV2	2936431	5296362	2.41%	0.616%
81	17.269	2425	2428	2432	rVB2	2643704	3170295	1.44%	0.369%
82	17.433	2450	2456	2462	rBV3	2911932	5526089	2.51%	0.643%
83	17.639	2486	2491	2497	rVB	6148702	8716901	3.96%	1.015%
84	17.722	2501	2505	2511	rVB	1810240	2790250	1.27%	0.325%
85	17.822	2519	2522	2526	rVB	2154055	2539943	1.15%	0.296%
86	18.069	2560	2564	2568	rVB	3197162	4316561	1.96%	0.502%
87	18.357	2608	2613	2619	rBV	4784835	6882256	3.13%	0.801%
88	19.016	2720	2725	2732	rVB4	4030290	7882562	3.58%	0.917%
89	19.598	2819	2824	2826	rBV	2709836	4032051	1.83%	0.469%
90	19.622	2826	2828	2833	rVB2	2654836	3001821	1.36%	0.349%
91	20.174	2917	2922	2926	rBV2	3399590	4696113	2.13%	0.547%
92	20.698	3006	3011	3015	rBV	2211763	2727846	1.24%	0.318%
93	21.216	3094	3099	3105	rVB	6856695	8971373	4.08%	1.044%
94	21.533	3148	3153	3160	rVB	1837781	2764581	1.26%	0.322%
95	21.780	3190	3195	3207	rVB	2283629	3512355	1.60%	0.409%
96	22.274	3273	3279	3286	rBV2	2506398	4034207	1.83%	0.470%
97	22.410	3295	3302	3308	rVB3	2336049	4634594	2.11%	0.539%
98	24.604	3667	3675	3688	rVB	1468752	4312579	1.96%	0.502%
99	24.827	3704	3713	3726	rVB2	1867296	5282751	2.40%	0.615%
100	27.474	4153	4163	4168	rBV4	681095	2330460	1.06%	0.271%

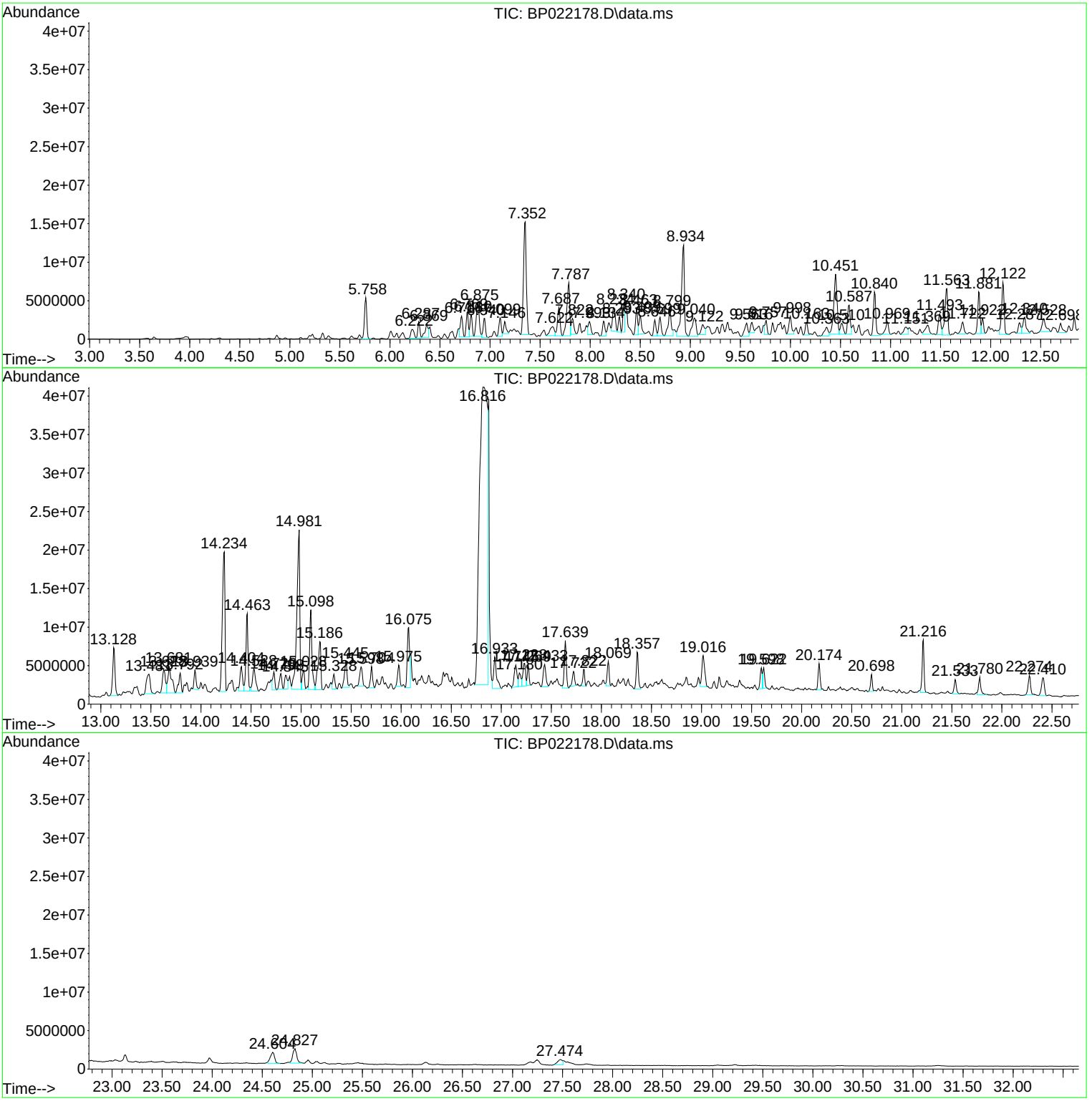
Sum of corrected areas: 859144702

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.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\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

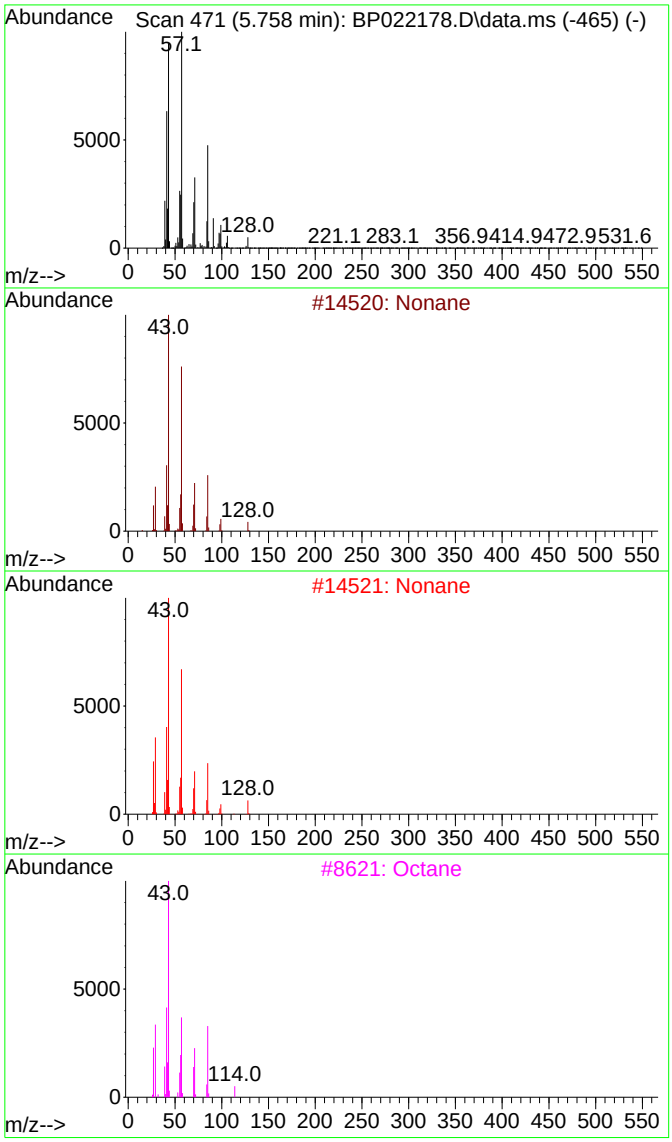
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

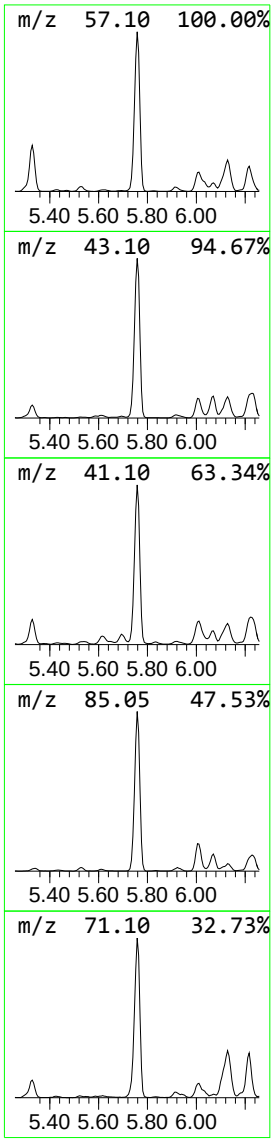
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.758	25.10 ng/ul	8399460	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane	128	C9H20	000111-84-2	95
2	Nonane	128	C9H20	000111-84-2	94
3	Octane	114	C8H18	000111-65-9	59
4	Octane	114	C8H18	000111-65-9	59
5	4,4-Dimethyl octane	142	C10H22	015869-95-1	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

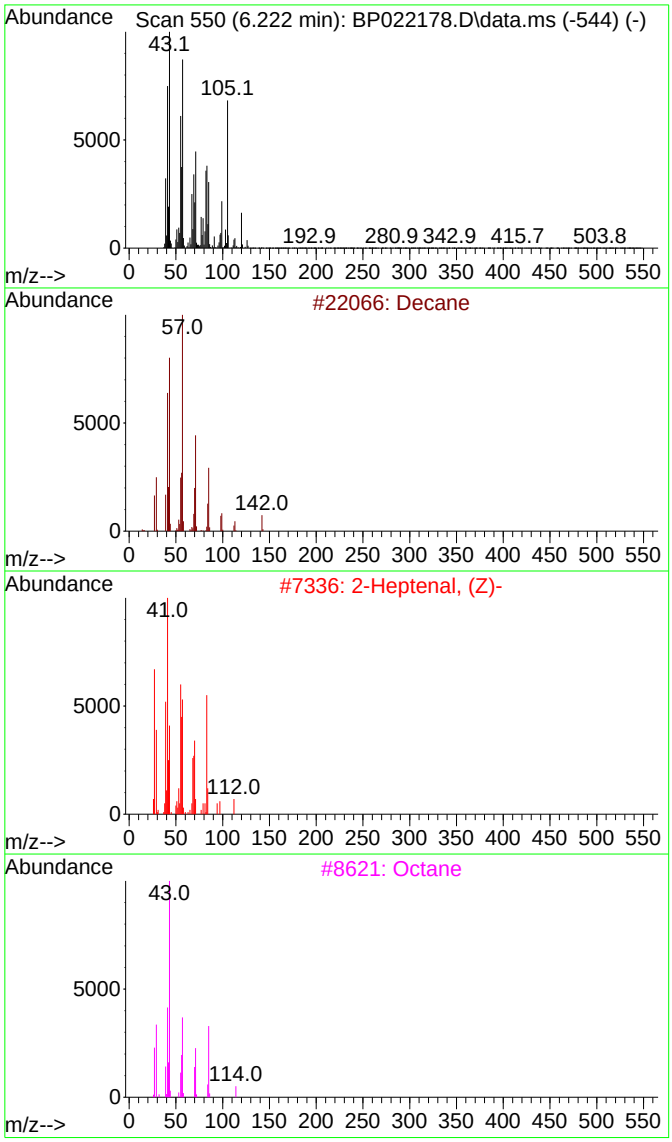
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

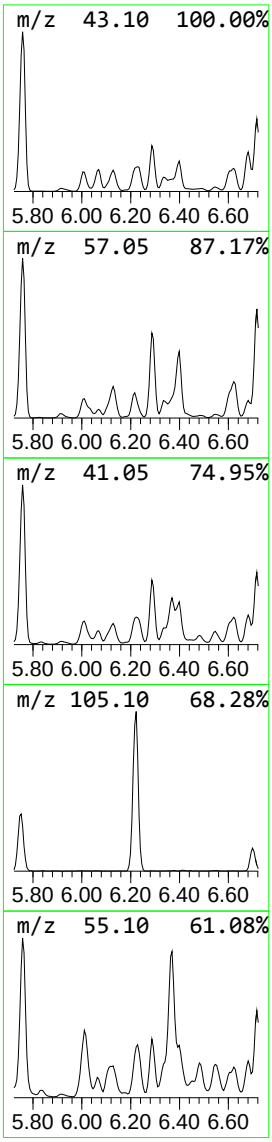
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.222	8.04 ng/ul	2689750	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	43
2	2-Heptenal, (Z)-	112	C7H12O	057266-86-1	38
3	Octane	114	C8H18	000111-65-9	38
4	Decane	142	C10H22	000124-18-5	38
5	Octane	114	C8H18	000111-65-9	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

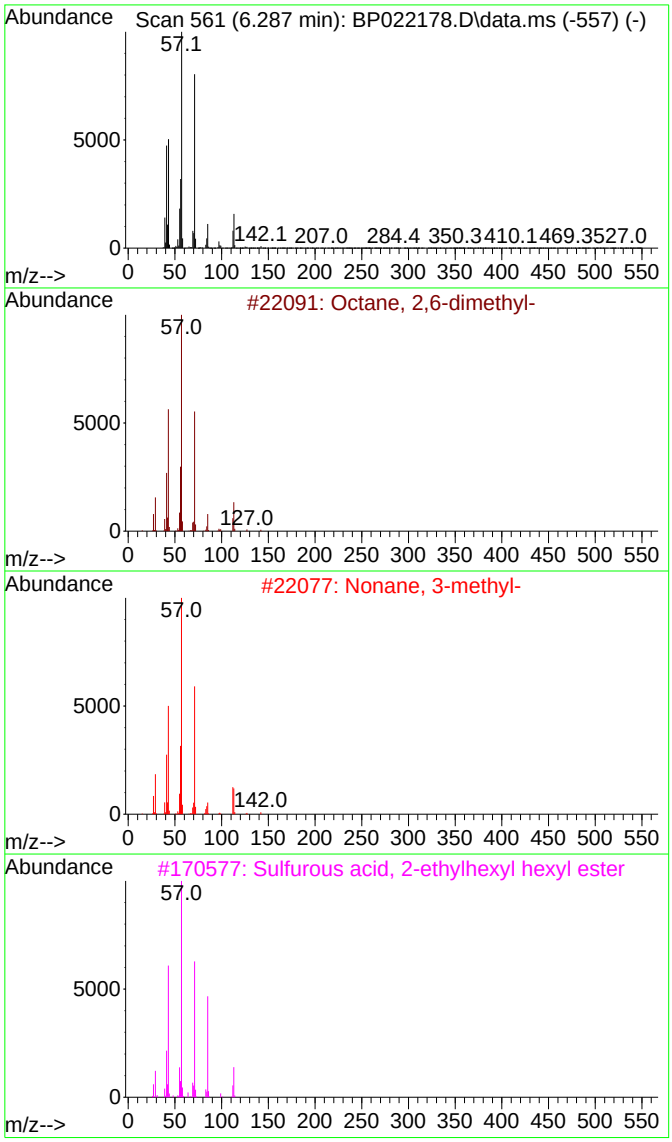
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

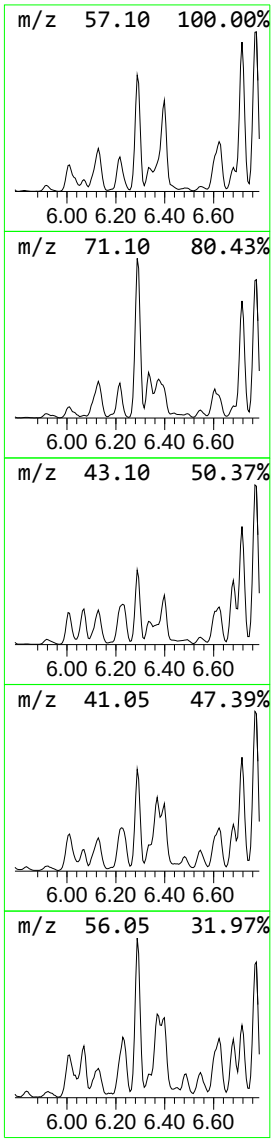
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.287	8.77 ng/ul	2934300	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	94
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	91
3	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	72
4	Octane, 3-ethyl-	142	C10H22	005881-17-4	72
5	Nonane, 3-methyl-	142	C10H22	005911-04-6	70





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

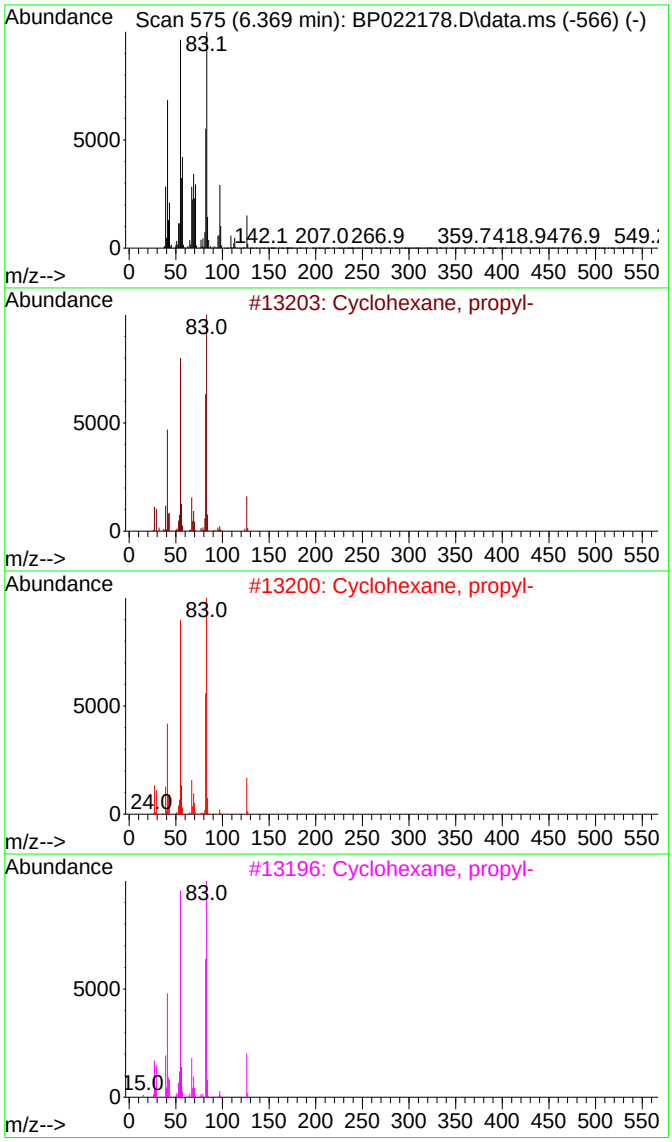
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

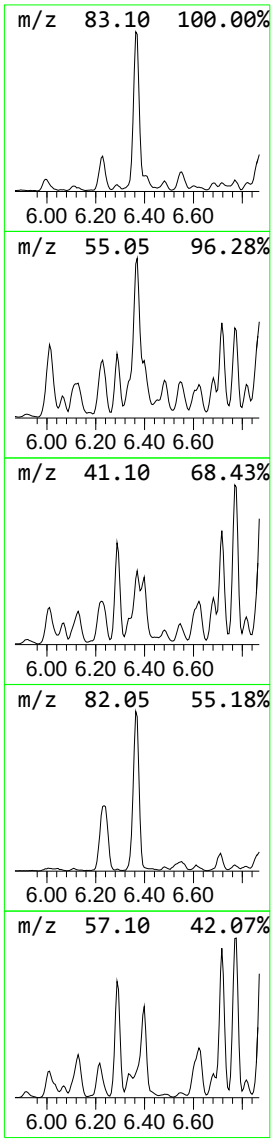
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Cyclic6.369 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.369	11.53 ng/ul	3858630	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	49
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	49
3	Cyclohexane, propyl-	126	C9H18	001678-92-8	49
4	Cyclohexane, propyl-	126	C9H18	001678-92-8	49
5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

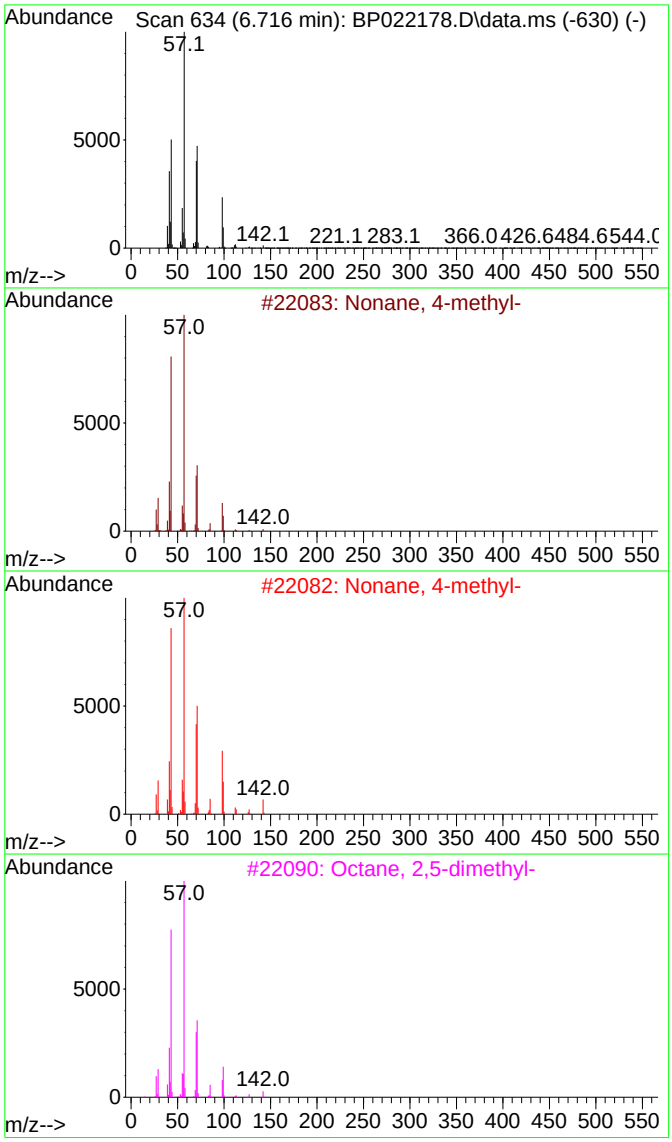
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

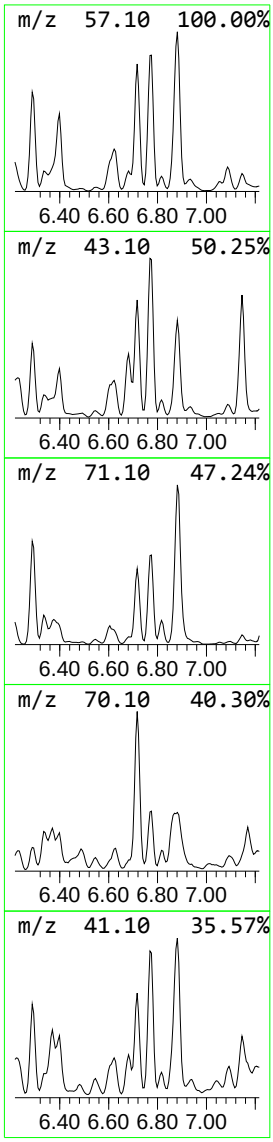
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.716	12.57 ng/ul	4206700	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-	142	C10H22	017301-94-9	72
2	Nonane, 4-methyl-	142	C10H22	017301-94-9	72
3	Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	64
4	Nonane, 4-methyl-	142	C10H22	017301-94-9	59
5	Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	59





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

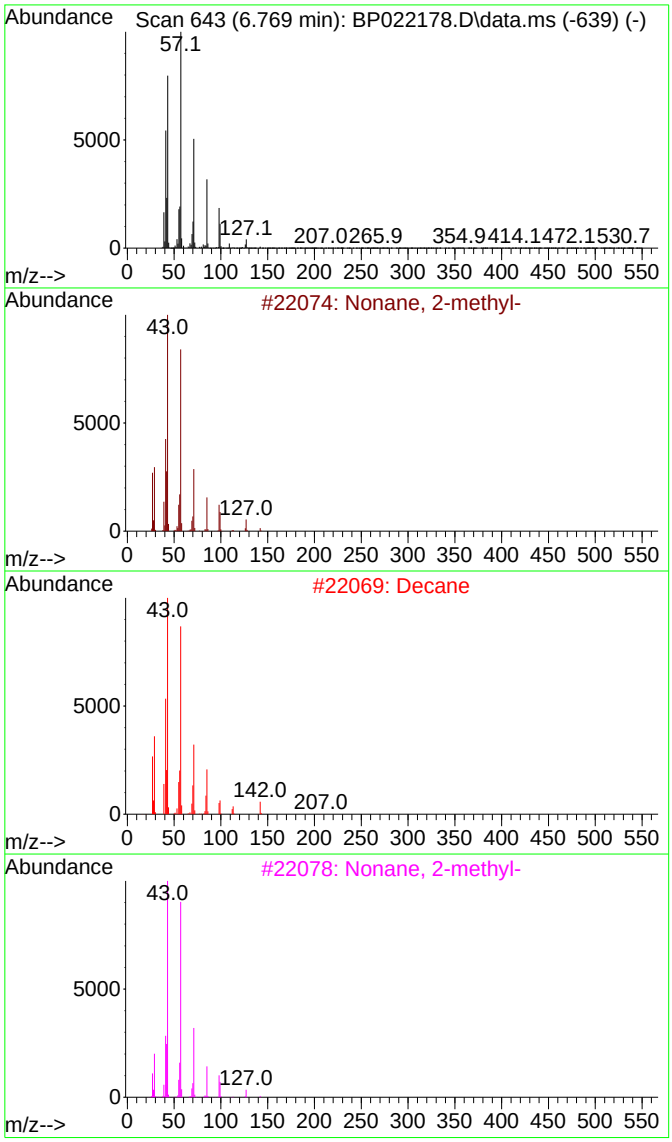
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

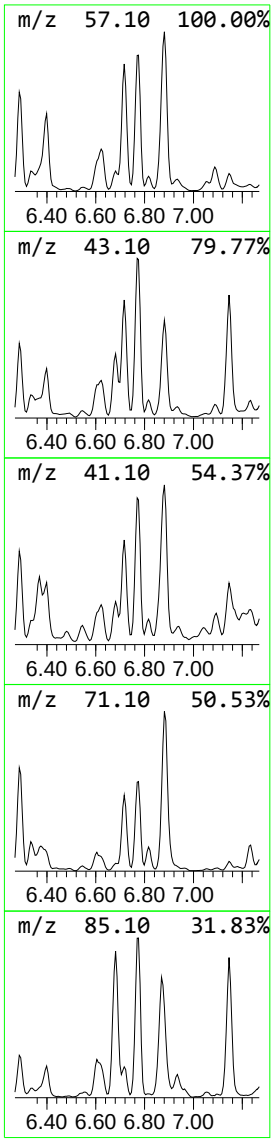
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.769	14.78 ng/ul	4945360	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2-methyl-	142	C10H22	000871-83-0	90
2	Decane	142	C10H22	000124-18-5	64
3	Nonane, 2-methyl-	142	C10H22	000871-83-0	58
4	Undecane, 3,7-dimethyl-	184	C13H28	017301-29-0	53
5	Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	53





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

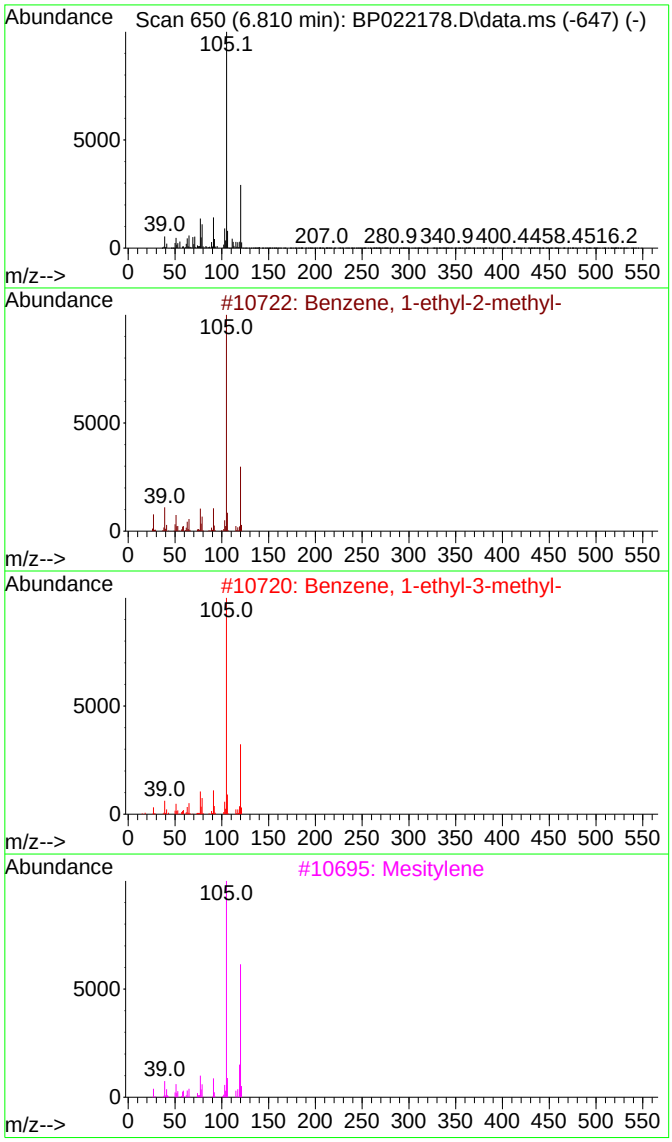
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

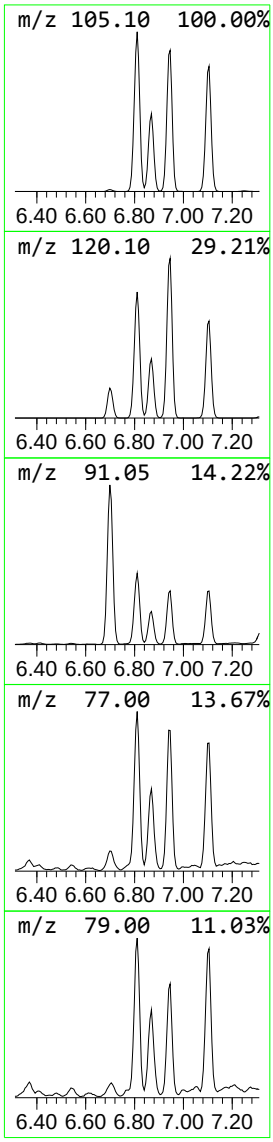
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.810	11.71 ng/ul	3919930	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

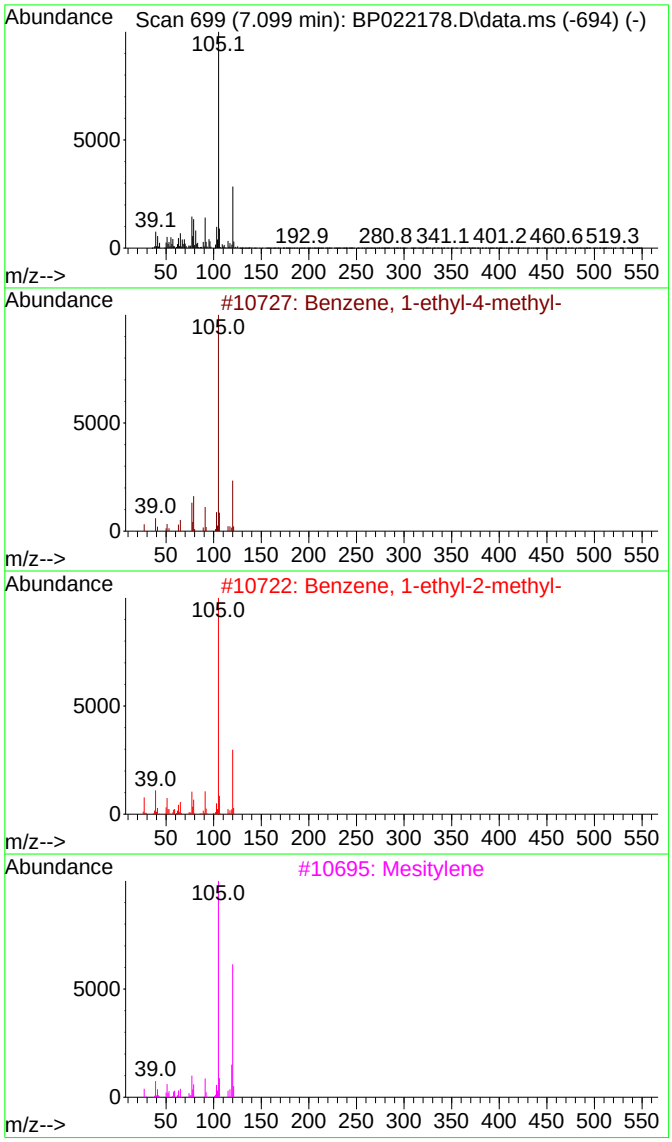
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

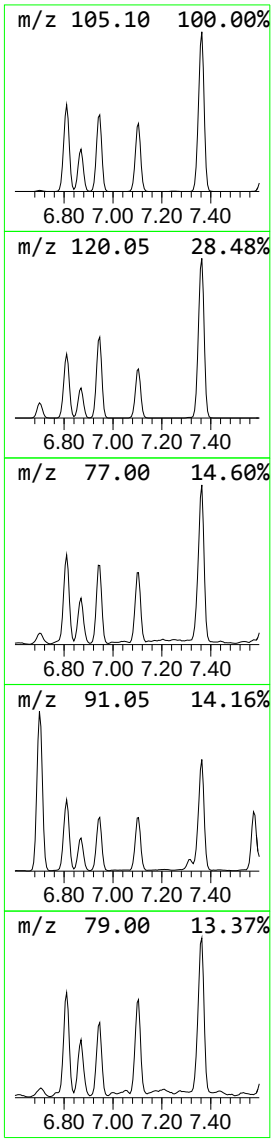
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.099	11.77 ng/ul	3940130	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

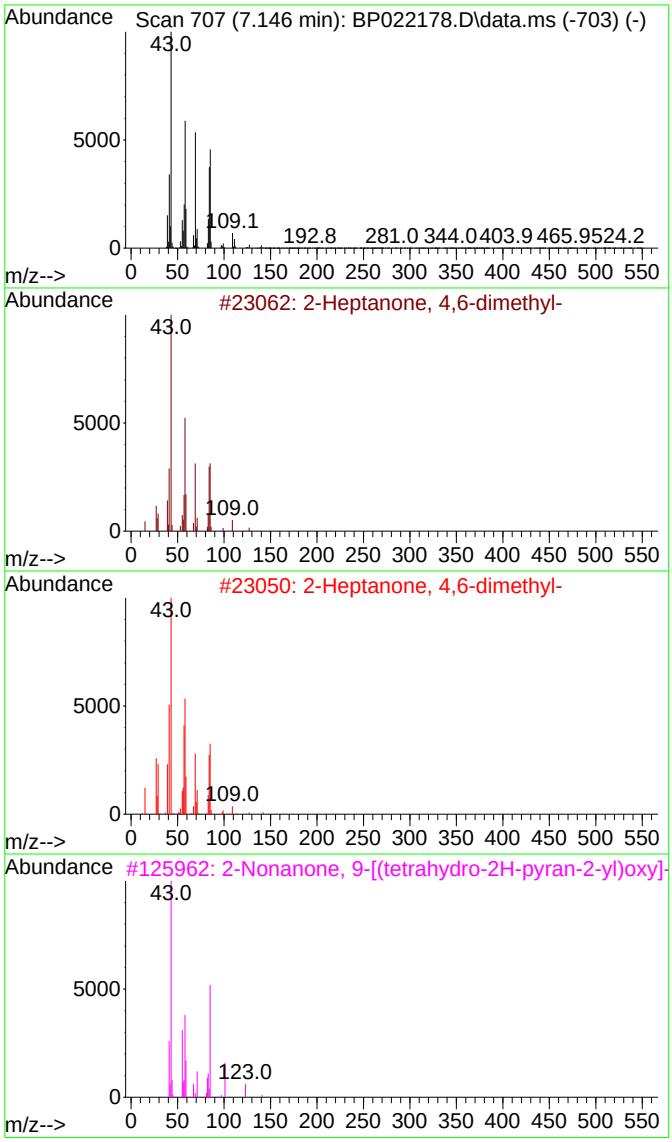
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

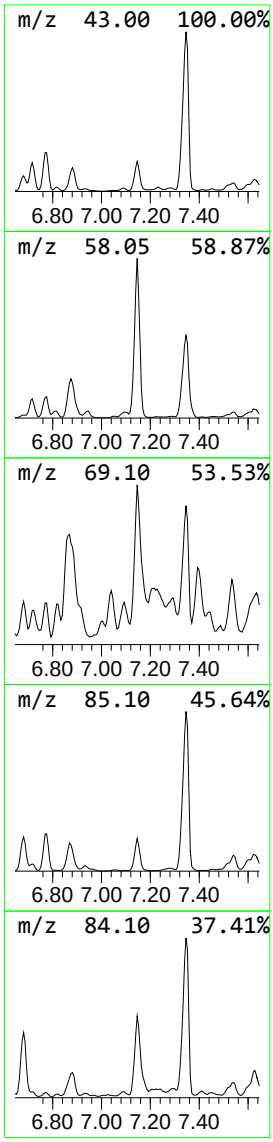
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2-Heptanone, 4,6-dimethyl- Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.146	7.76 ng/ul	2595520	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	80
2	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	72
3	2-Nonanone, 9-[(tetrahydro-2H-py...	242	C14H26O3	054699-41-1	64
4	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	47
5	Carbonic acid, hexyl prop-1-en-2...	186	C10H18O3	1000382-54-2	43





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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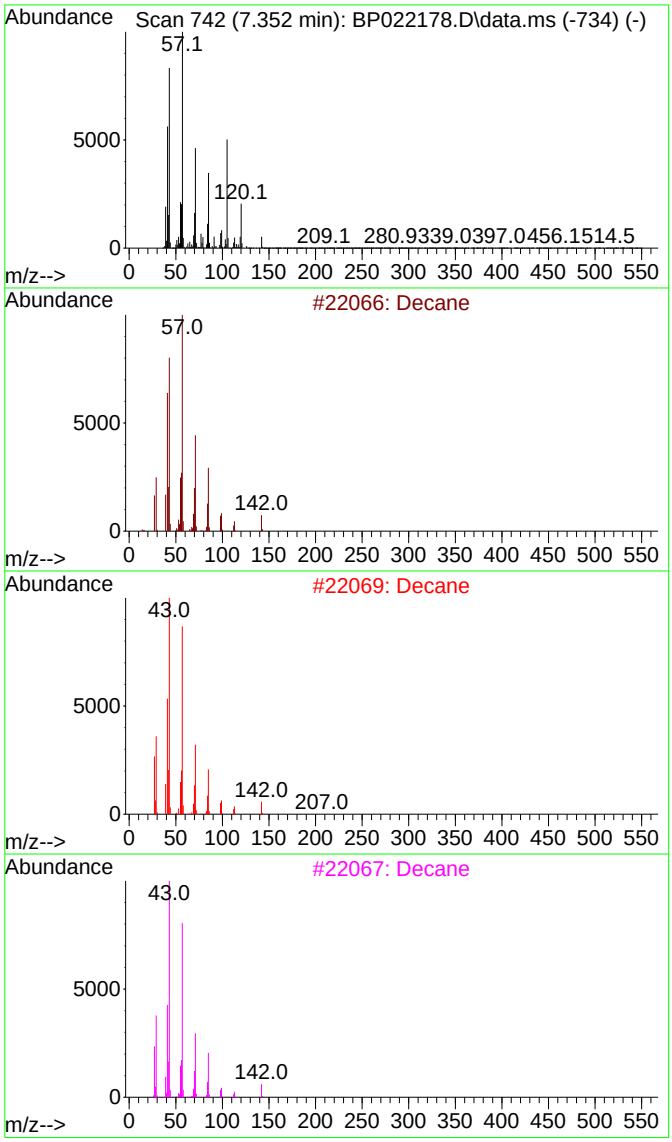
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

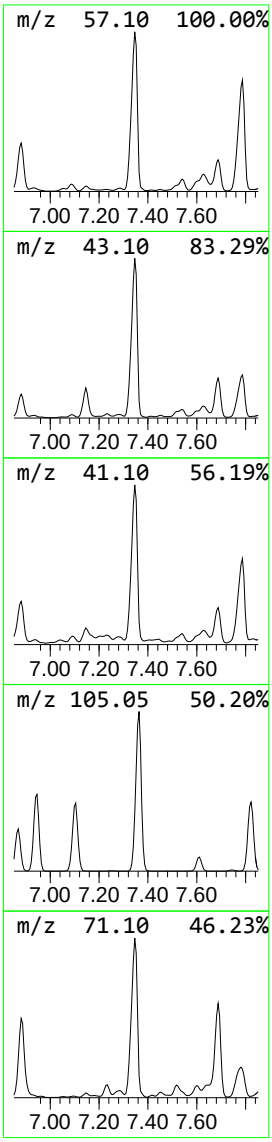
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Carbonic acid, prop-1-en-2-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.352	84.49 ng/ul	28272900	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	94
2	Decane	142	C10H22	000124-18-5	93
3	Decane	142	C10H22	000124-18-5	93
4	Decane	142	C10H22	000124-18-5	81
5	Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	53





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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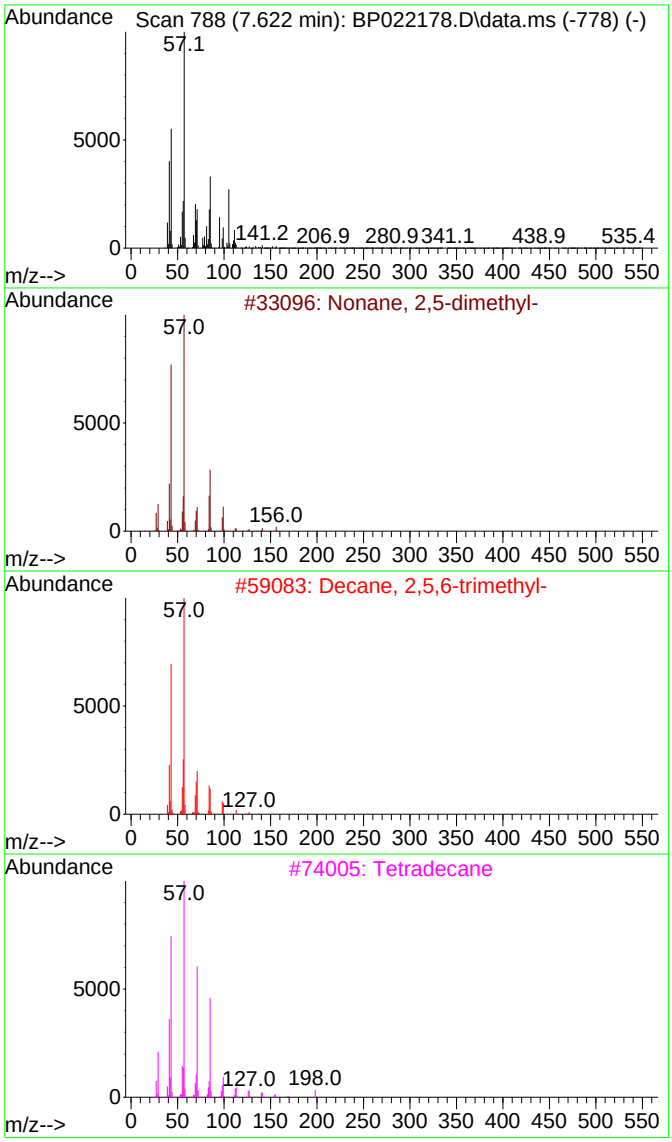
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

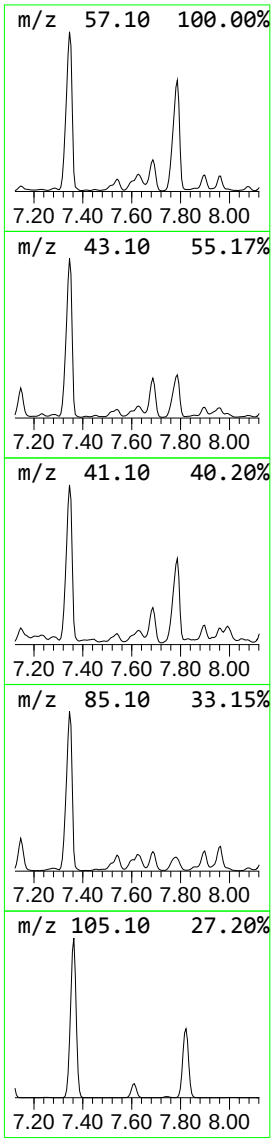
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.622	10.95 ng/ul	3663830	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	47
2	Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	43
3	Tetradecane	198	C14H30	000629-59-4	38
4	Pentane, 3-ethyl-2,3-dimethyl-	128	C9H20	016747-33-4	38
5	Tridecane	184	C13H28	000629-50-5	38





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

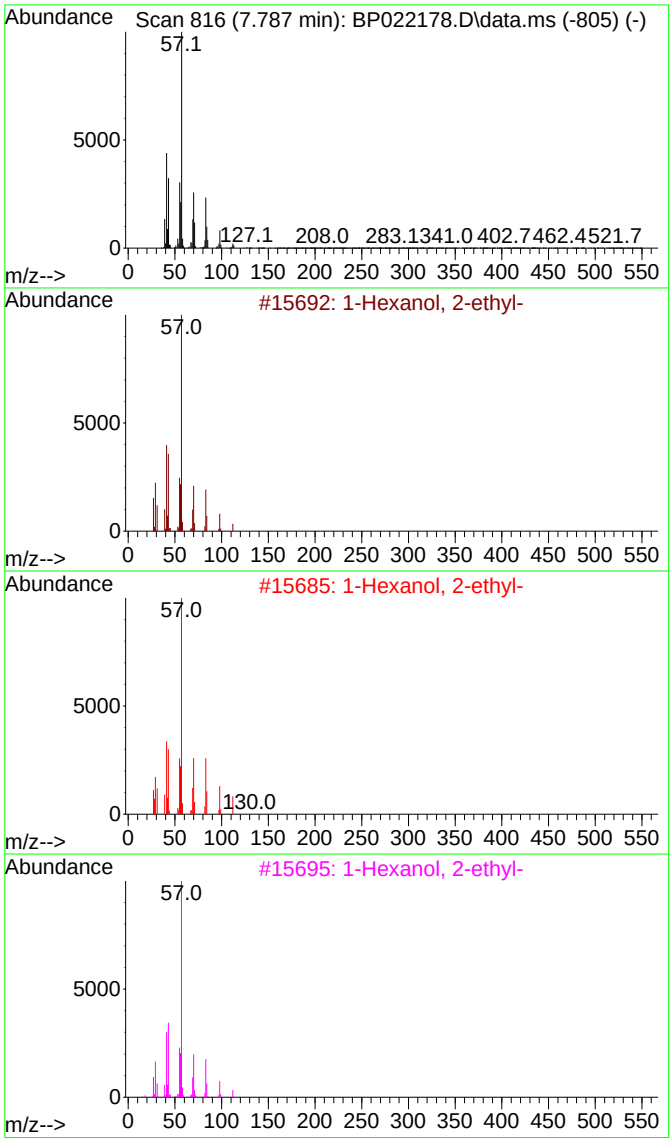
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

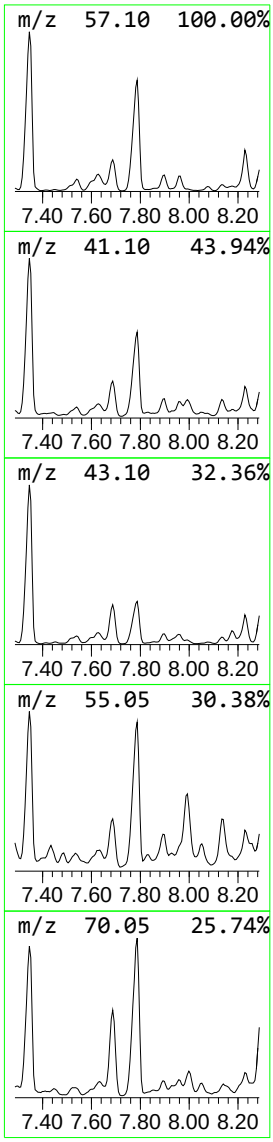
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.787	42.20 ng/ul	14121600	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
5	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53





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Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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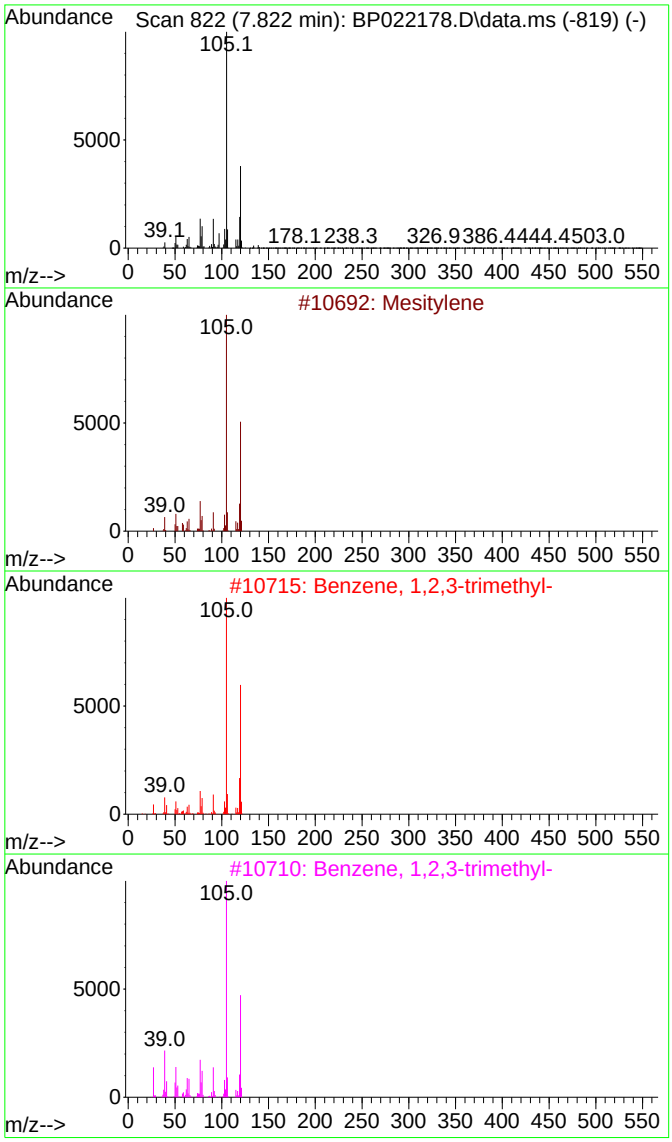
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

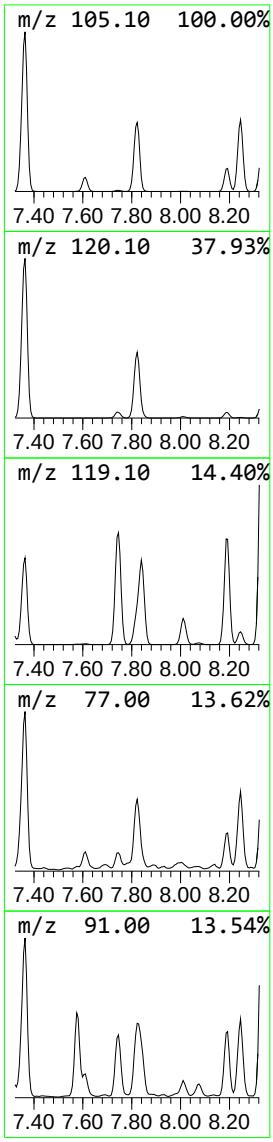
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Mesitylene Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.822	10.83 ng/ul	3625450	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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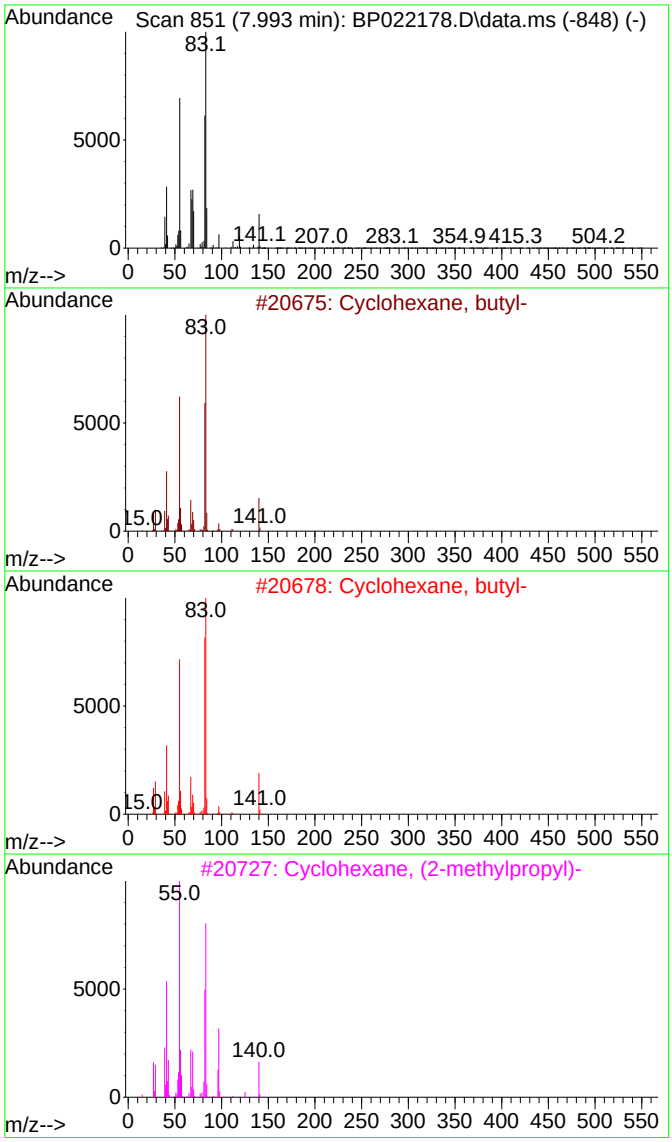
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

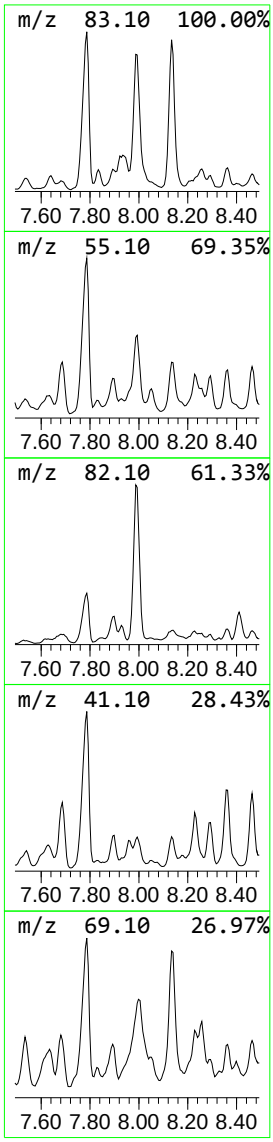
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 (DEL) Alkane: Cyclic7.993 Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.993	8.76 ng/ul	2930650	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, butyl-	140	C10H20	001678-93-9	74
2	Cyclohexane, butyl-	140	C10H20	001678-93-9	64
3	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	59
4	Cyclohexane, octyl-	196	C14H28	001795-15-9	59
5	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	53





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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

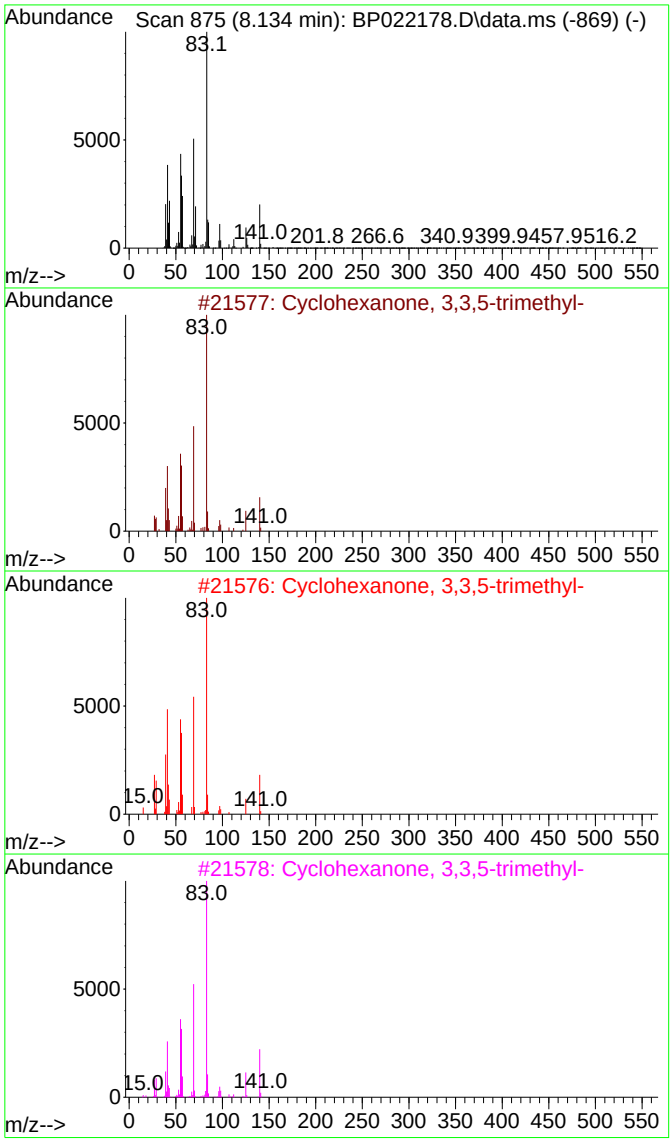
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

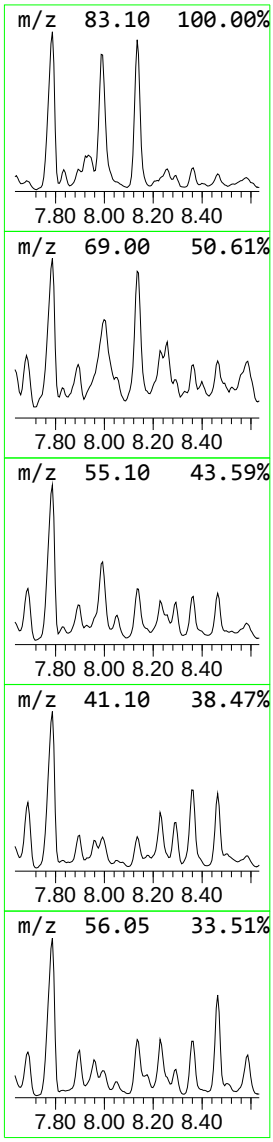
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Cyclohexanone, 3,3,5-trimet... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.134	9.97 ng/ul	3335080	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	76
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	76
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	59
5	1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	52





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Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

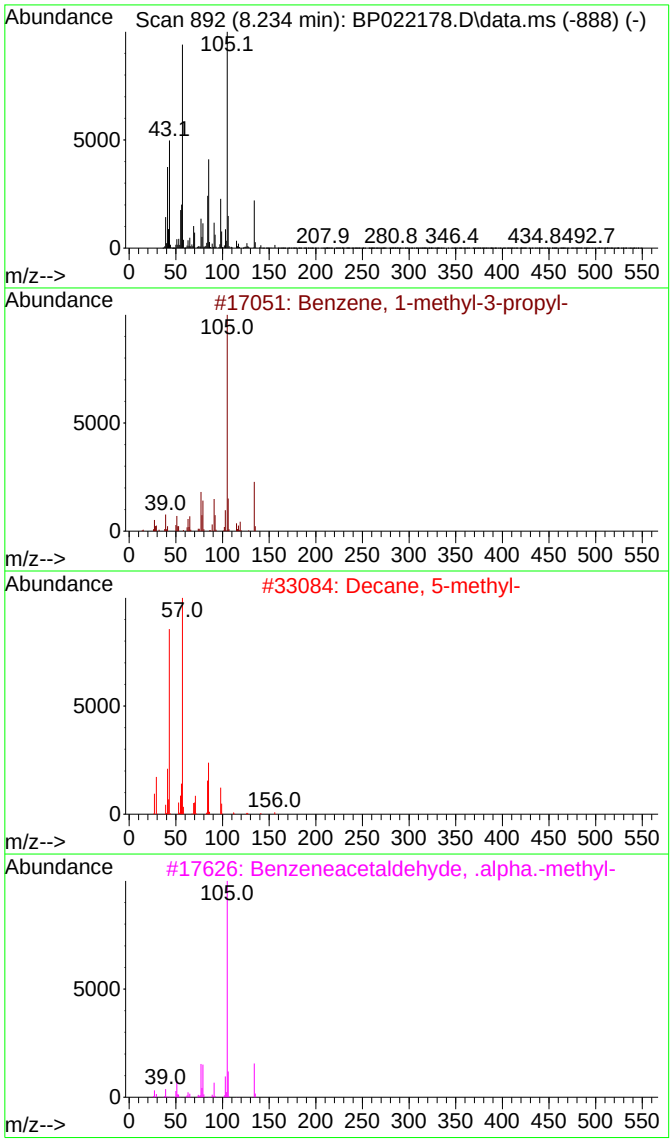
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

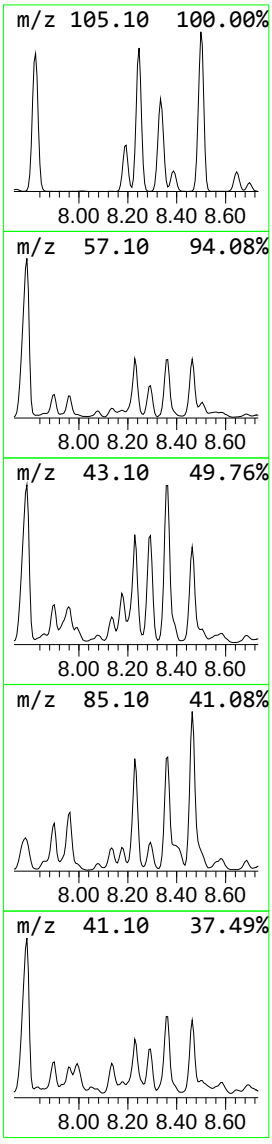
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 1-methyl-3-propyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	17.24 ng/ul	5768090	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	64
2	Decane, 5-methyl-	156	C11H24	013151-35-4	46
3	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	35
4	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	35
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	35





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

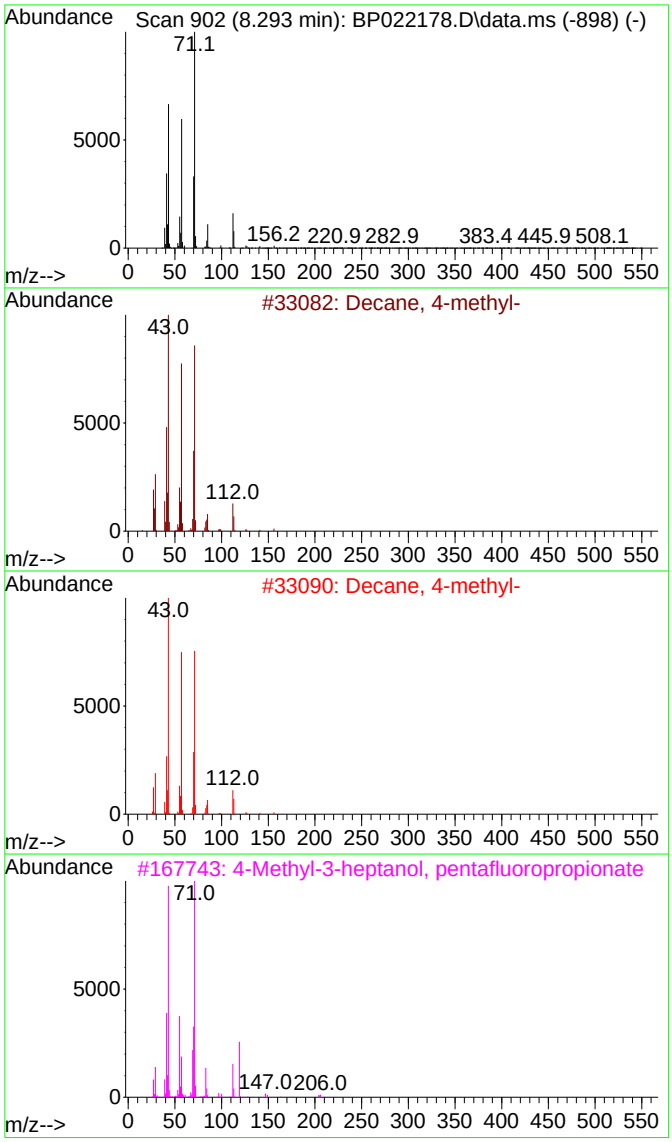
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

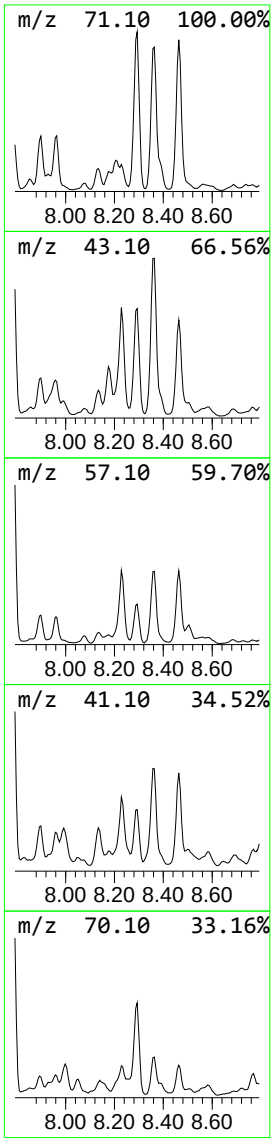
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.293	7.69 ng/ul	2574630	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-	156	C11H24	002847-72-5	91
2	Decane, 4-methyl-	156	C11H24	002847-72-5	87
3	4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	72
4	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72
5	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

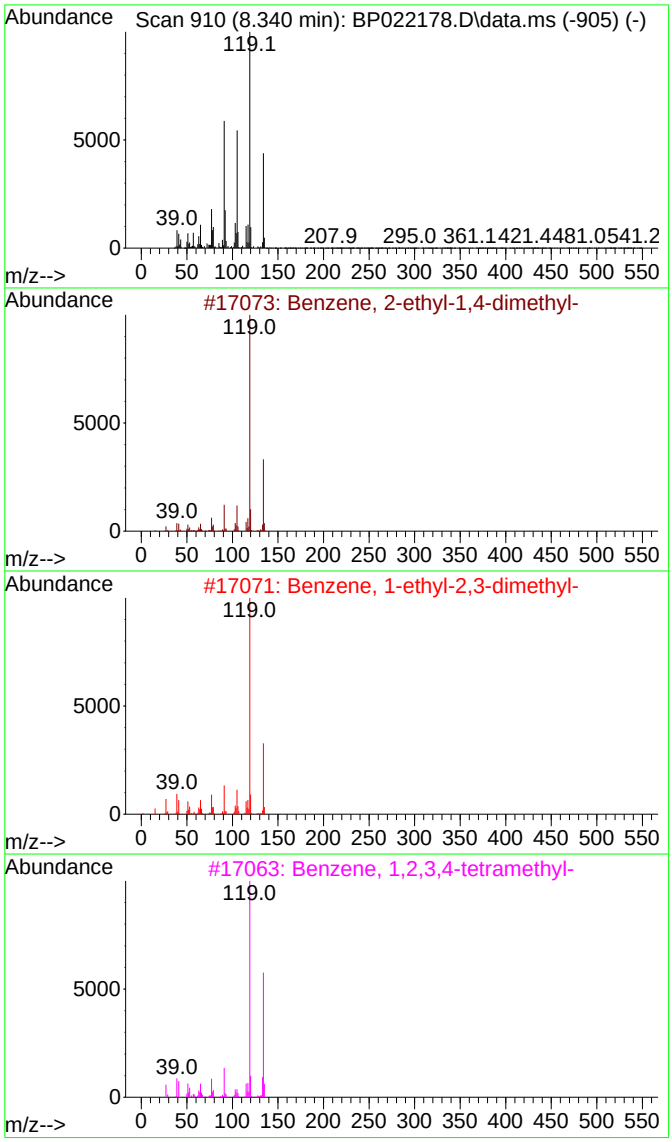
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

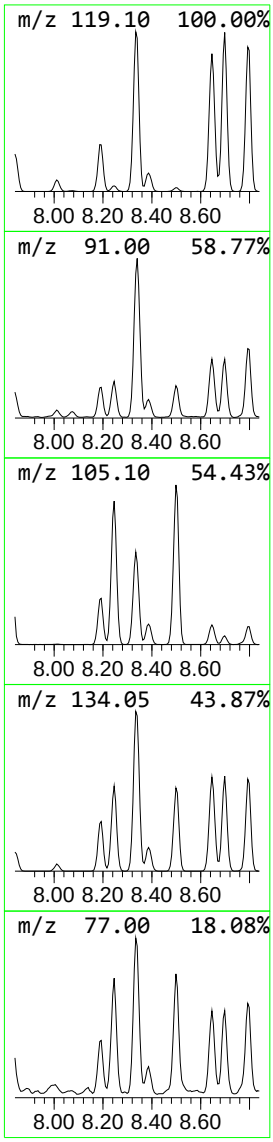
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.340	15.76 ng/ul	5273580	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

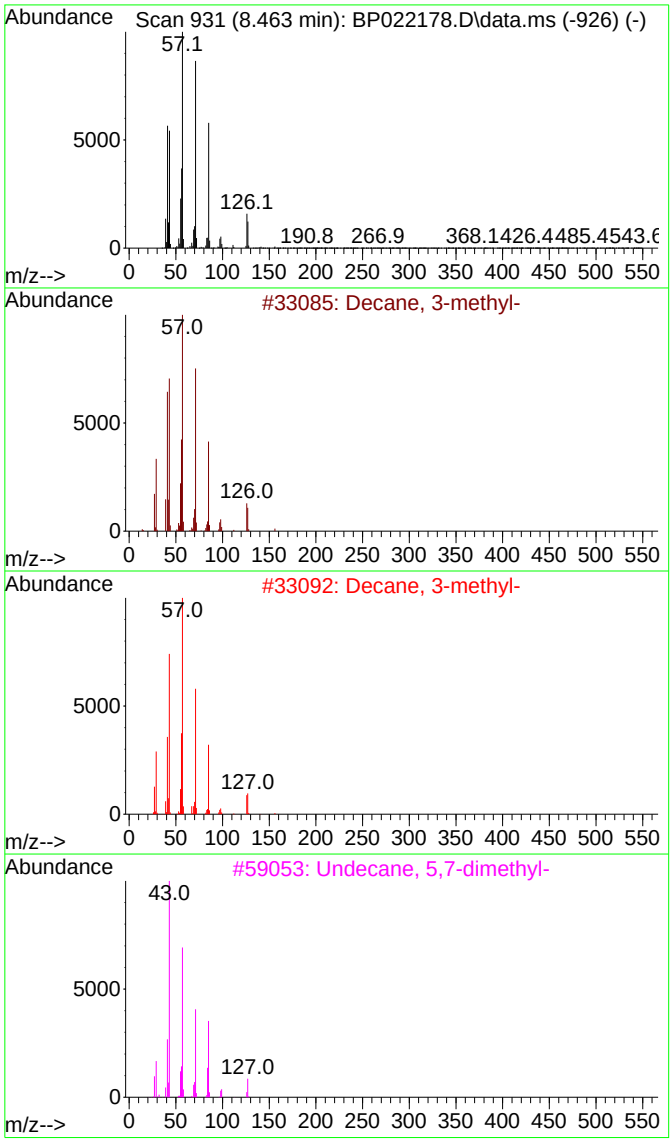
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

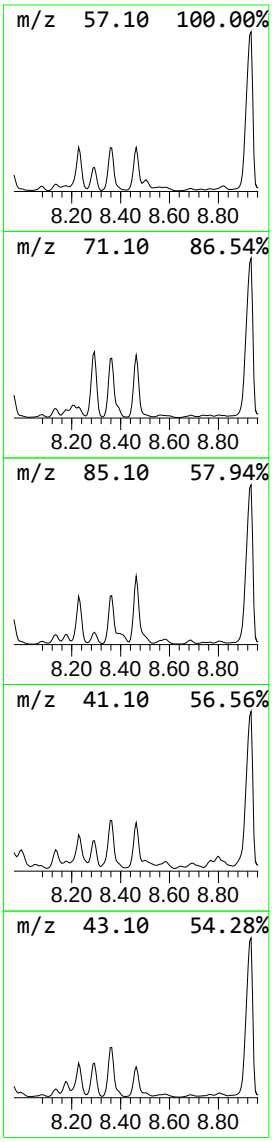
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.463	15.05 ng/ul	5035070	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3-methyl-	156	C11H24	013151-34-3	95
2	Decane, 3-methyl-	156	C11H24	013151-34-3	91
3	Undecane, 5,7-dimethyl-	184	C13H28	017312-83-3	78
4	Nonane, 5-butyl-	184	C13H28	017312-63-9	72
5	Nonane, 1-iodo-	254	C9H19I	004282-42-2	64





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

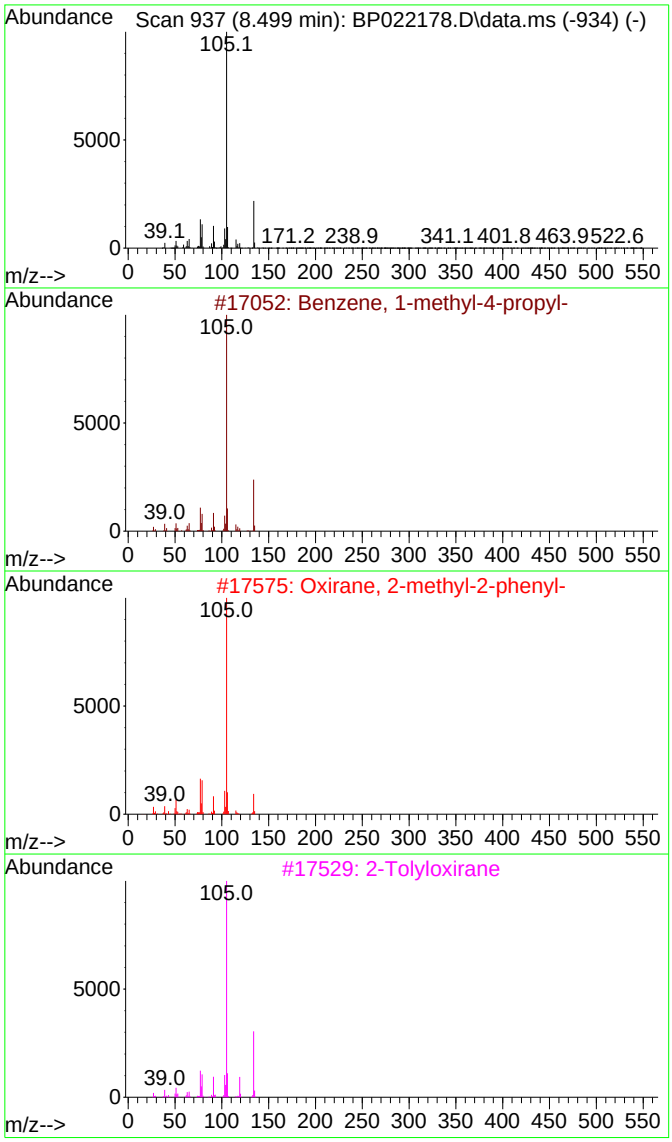
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

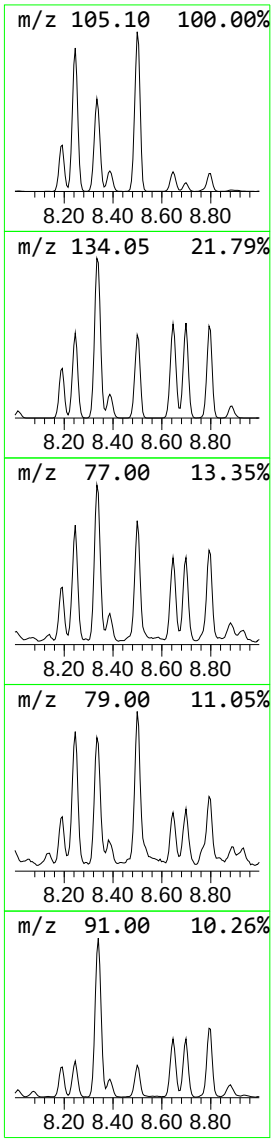
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 1-methyl-4-propyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.499	11.73 ng/ul	3924670	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2	Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	91
3	2-Tolyloxirane	134	C9H10O	002783-26-8	91
4	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	91
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

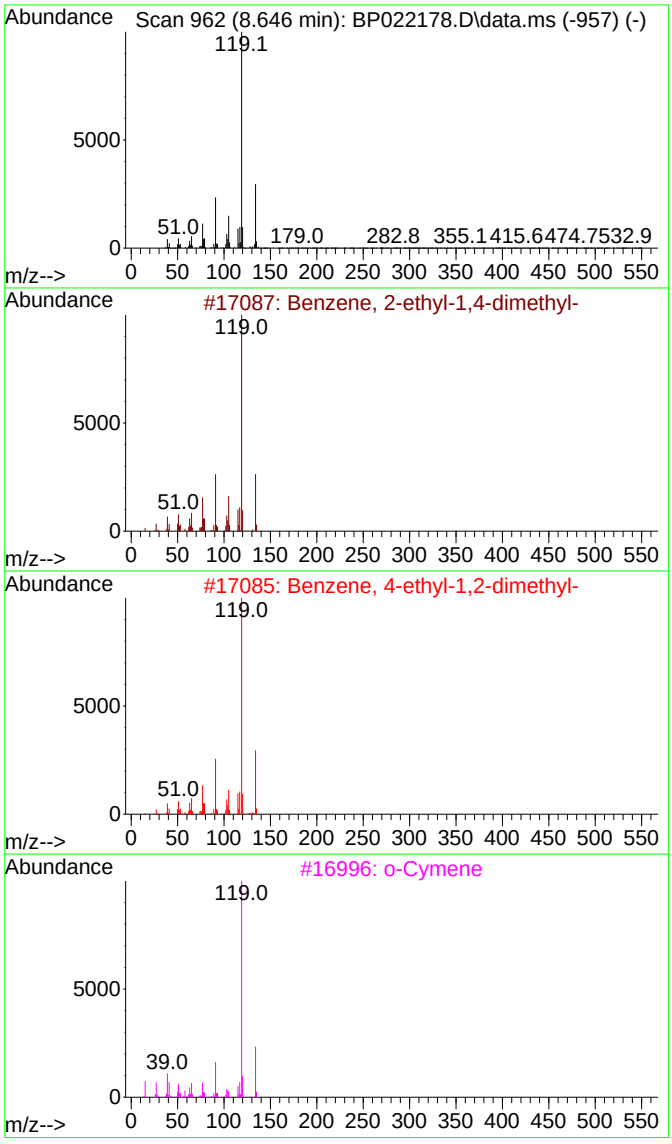
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

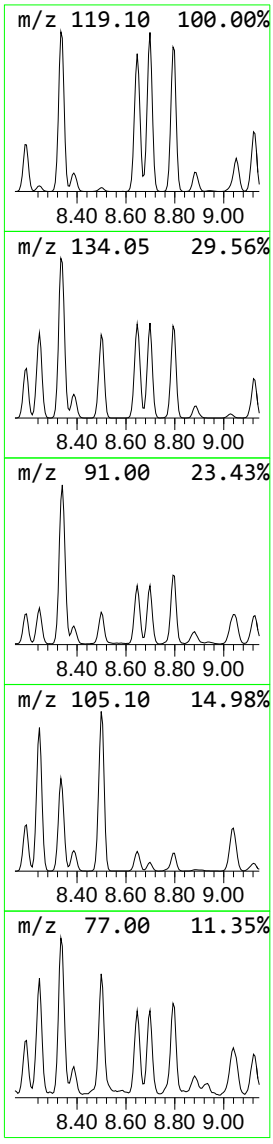
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.646	9.80 ng/ul	3280260	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	o-Cymene	134	C10H14	000527-84-4	95





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

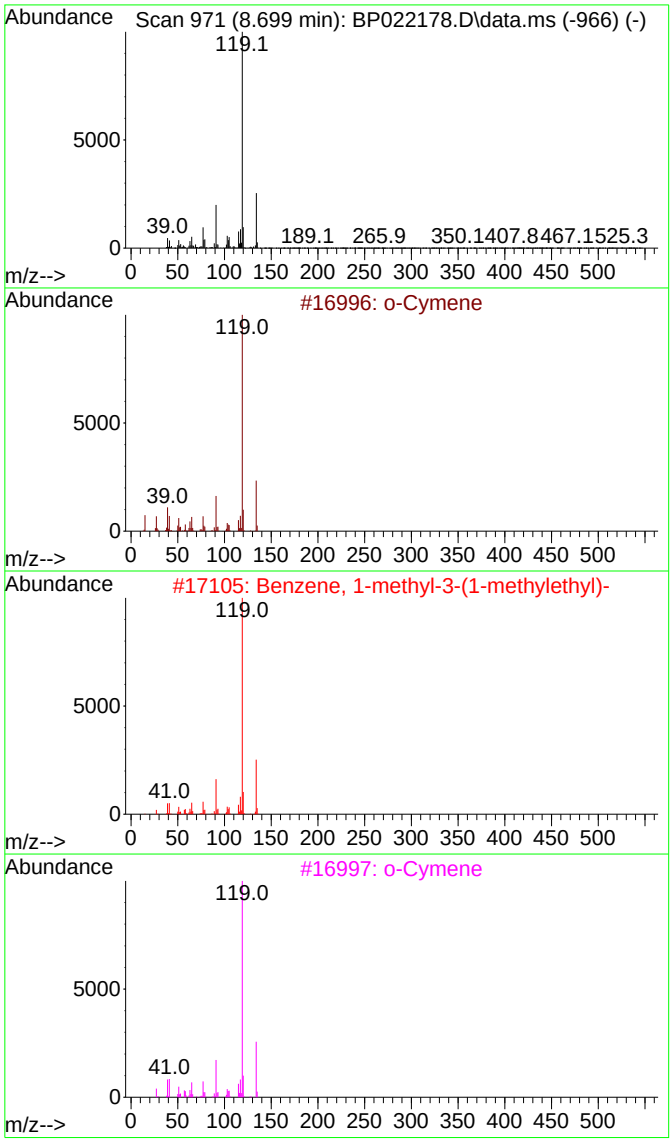
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

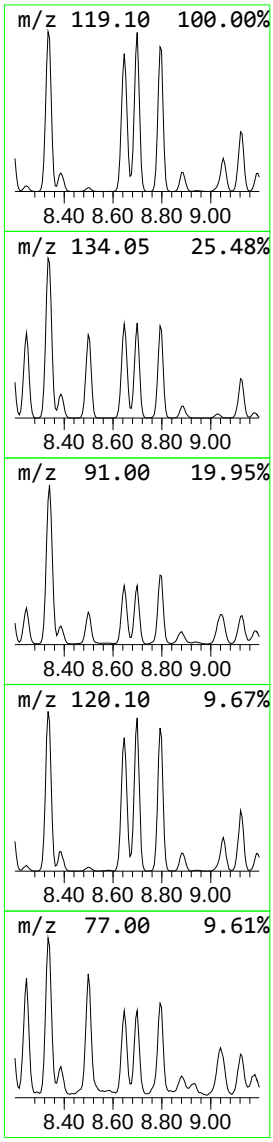
****

Peak Number 22 o-Cymene Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.699	12.01 ng/ul	4019810	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
3	o-Cymene	134	C10H14	000527-84-4	97
4	p-Cymene	134	C10H14	000099-87-6	97
5	o-Cymene	134	C10H14	000527-84-4	95





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

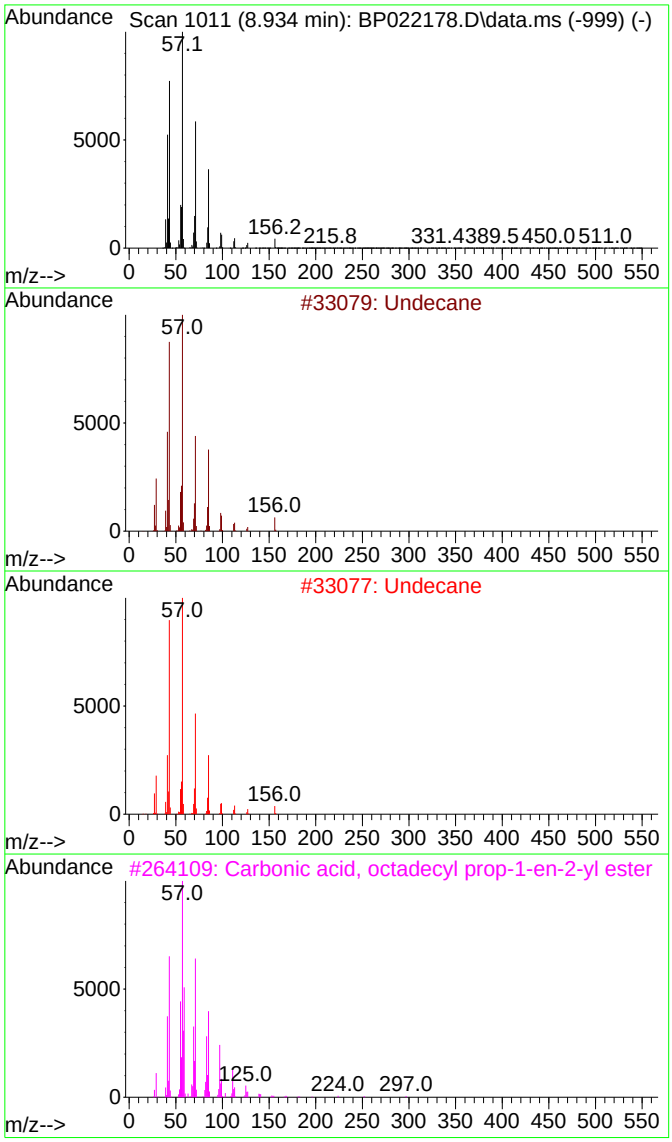
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

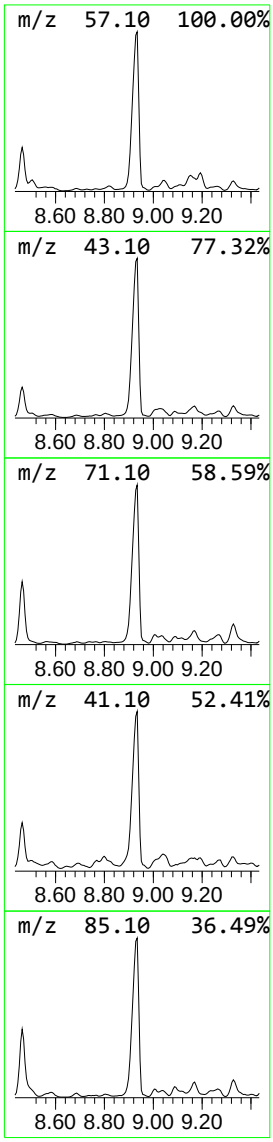
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.934	72.95 ng/ul	24411600	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	94
2	Undecane	156	C11H24	001120-21-4	91
3	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	90
4	Hexadecane	226	C16H34	000544-76-3	90
5	Tetradecane	198	C14H30	000629-59-4	90





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

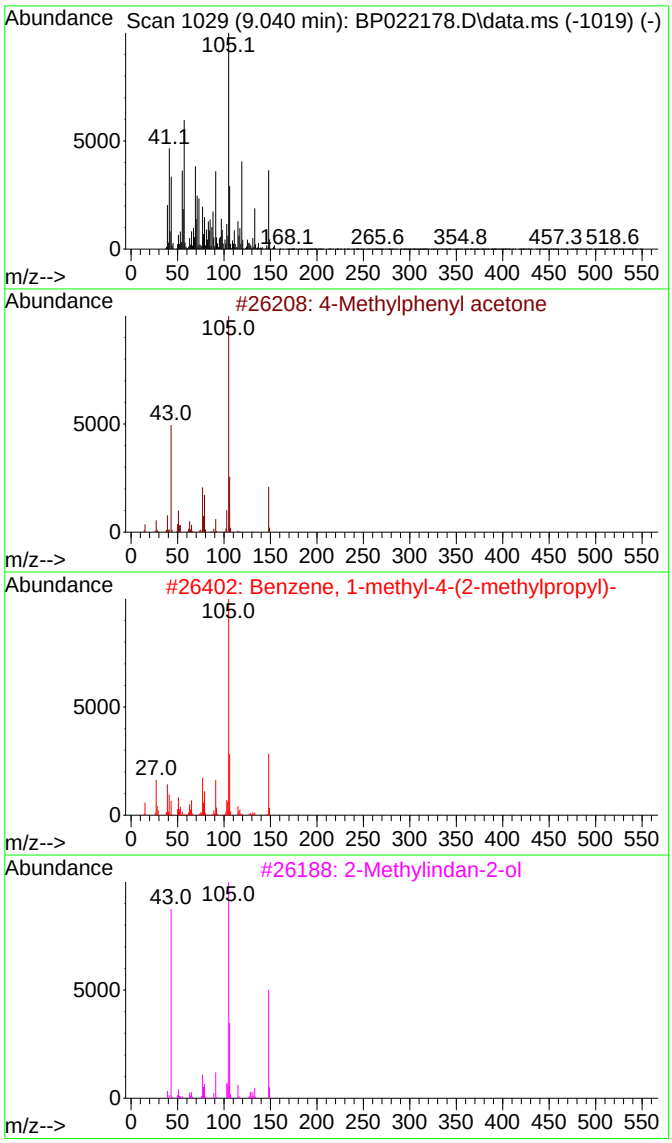
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

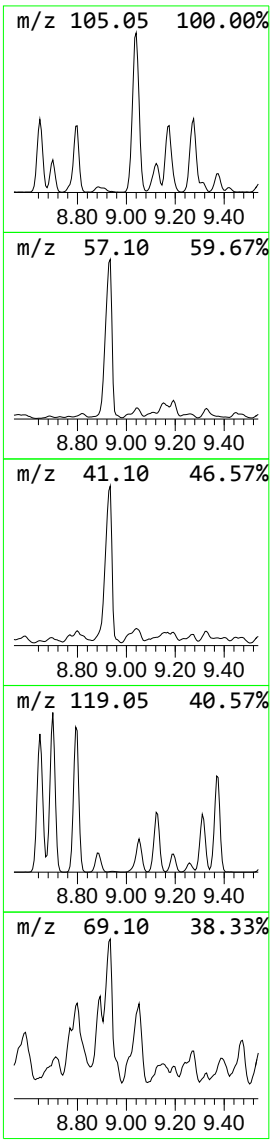
****

Peak Number 24 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.040	20.75 ng/ul	6944180	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methylphenyl acetone	148	C10H12O	002096-86-8	42
2	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	41
3	2-Methylindan-2-ol	148	C10H12O	033223-84-6	27
4	Benzene, 1-methyl-4-butyl	148	C11H16	001595-05-7	25
5	3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	25





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

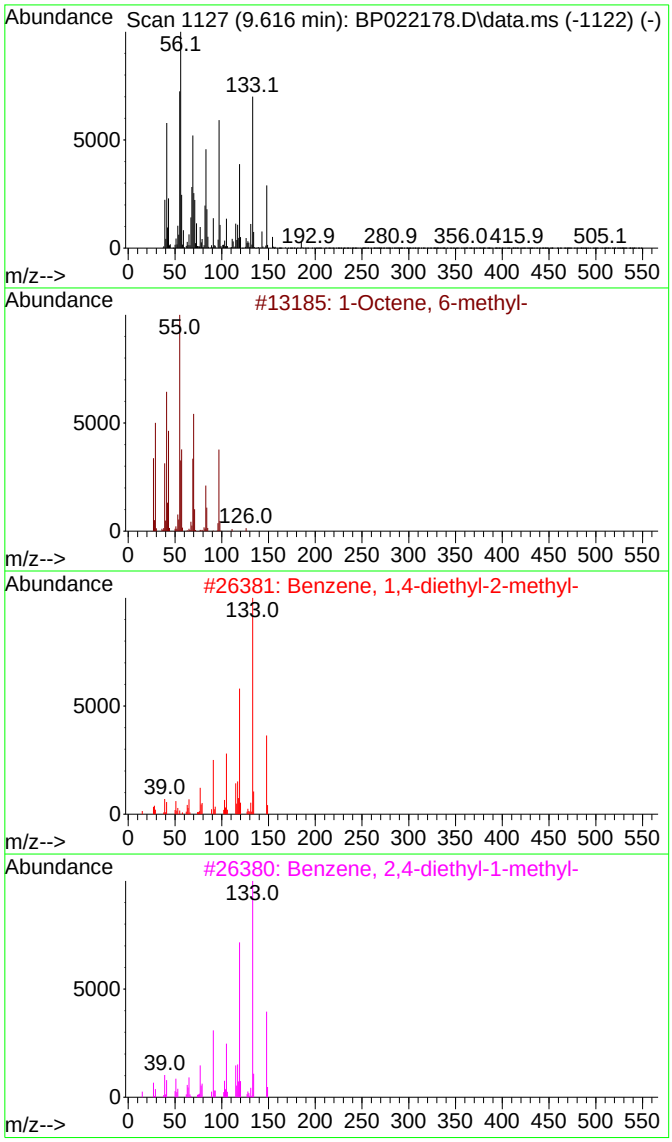
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

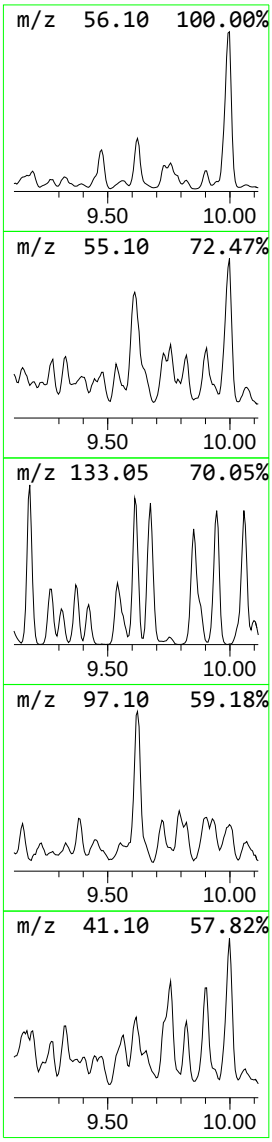
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.616	2.68 ng/ul	2390280	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octene, 6-methyl-	126	C9H18	013151-10-5	38
2	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	38
3	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	38
4	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	38
5	Cycloheptane	98	C7H14	000291-64-5	38





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

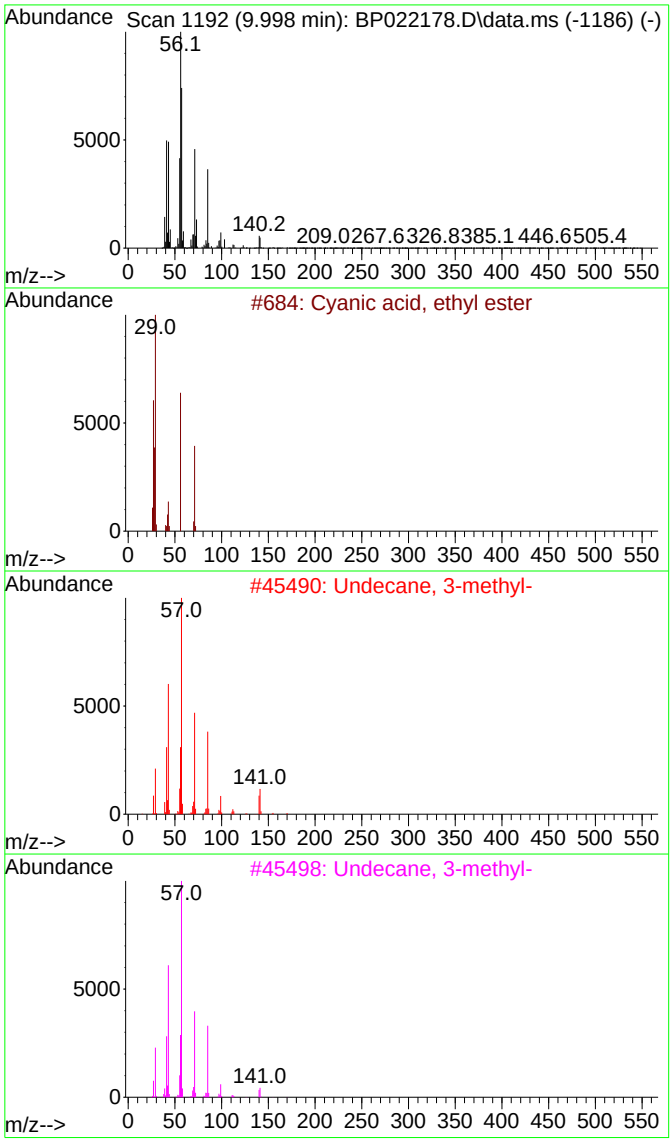
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

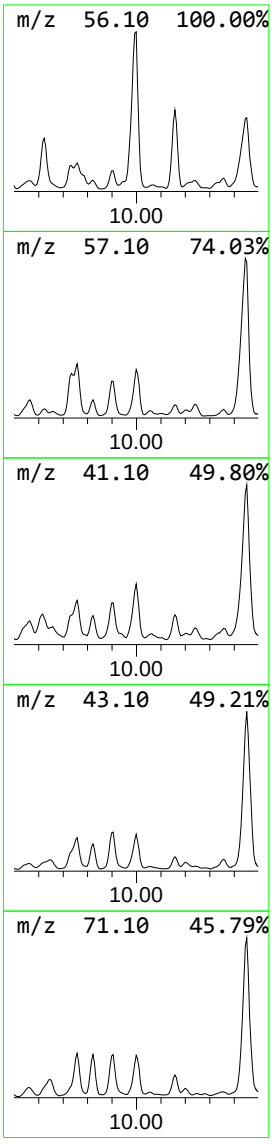
****

Peak Number 28 unknown-02 Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.998	4.79 ng/ul	4276640	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	49
2	Undecane, 3-methyl-	170	C12H26	001002-43-3	46
3	Undecane, 3-methyl-	170	C12H26	001002-43-3	43
4	1-Hexanol, 2-ethyl-2-propyl-	172	C11H24O	054461-00-6	38
5	2,4-Dimethyl-1-hexene	112	C8H16	016746-87-5	38





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

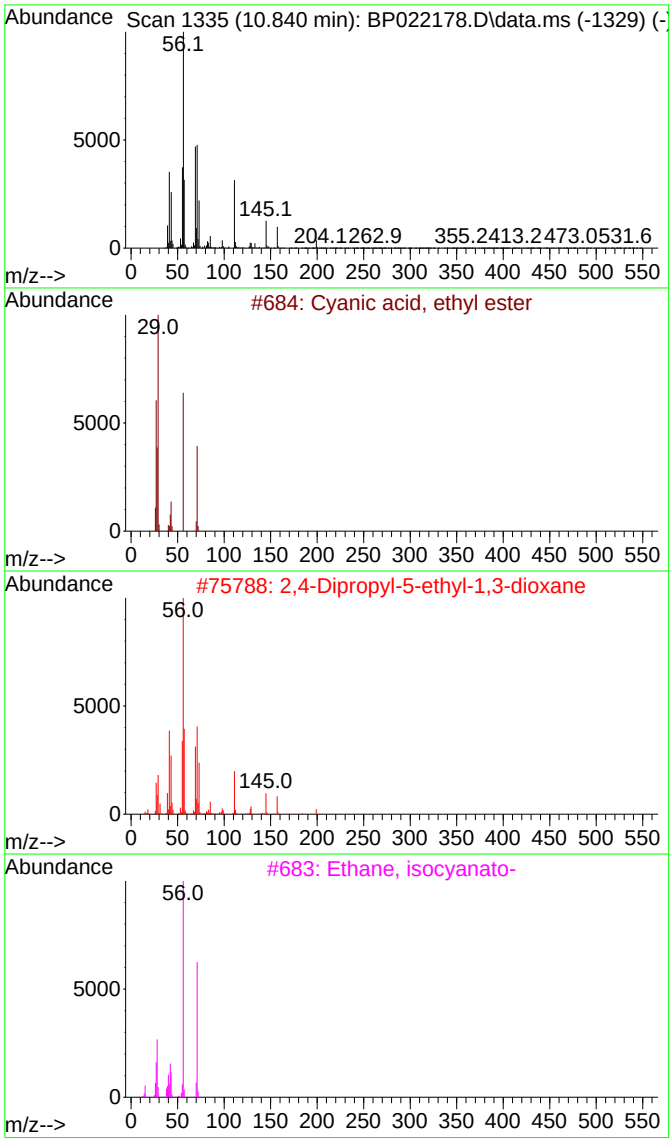
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

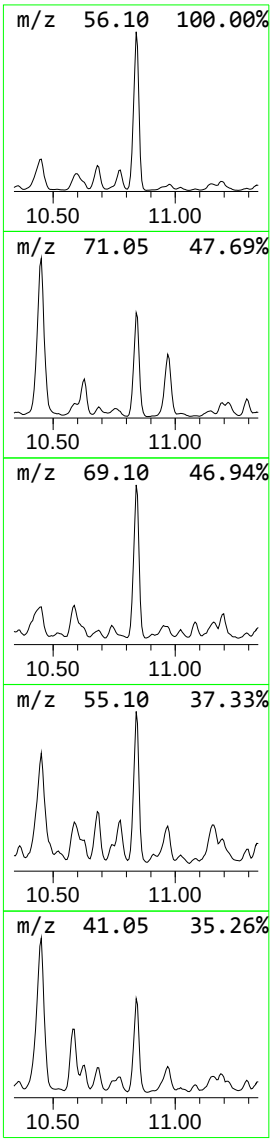
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 unknown-03 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.840	10.45 ng/ul	9324640	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	49
2	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	45
3	Ethane, isocyanato-	71	C3H5NO	000109-90-0	38
4	Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	32
5	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

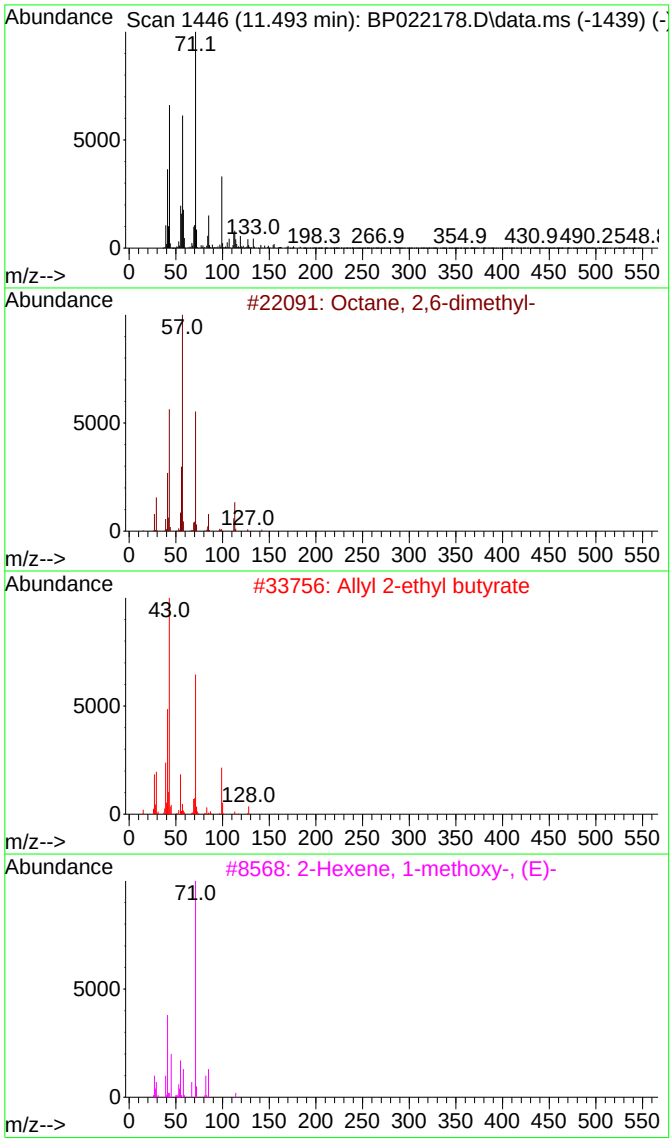
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

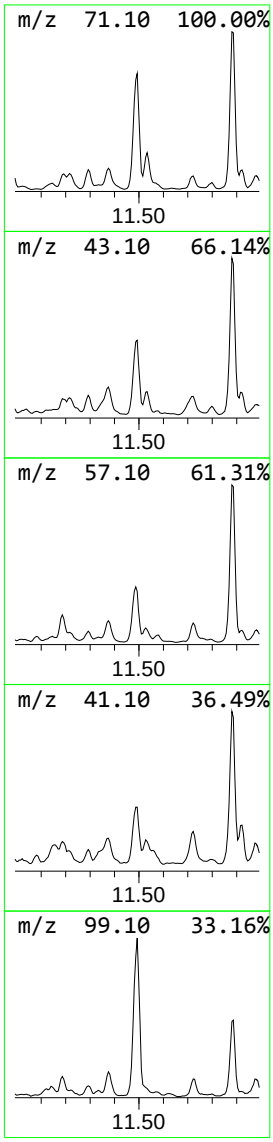
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.493	5.73 ng/ul	5113300	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	58
2	Allyl 2-ethyl butyrate	156	C9H16O2	007493-69-8	53
3	2-Hexene, 1-methoxy-, (E)-	114	C7H14O	056052-83-6	52
4	2-Hexene, 1-methoxy-, (E)-	114	C7H14O	056052-83-6	50
5	4-Heptanone, 3-ethyl-	142	C9H18O	001528-25-2	50





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

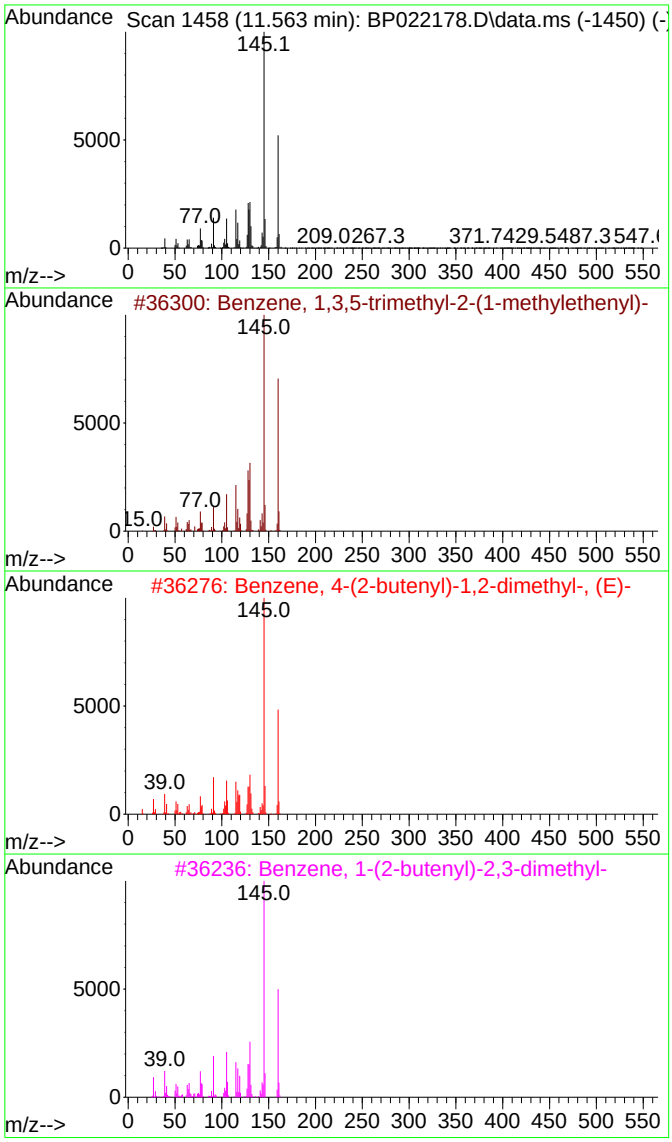
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

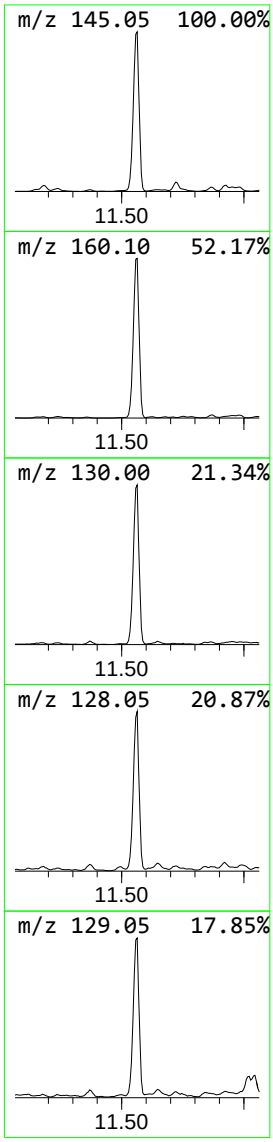
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 35 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.563	12.96 ng/ul	11561300	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	96
2	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	94
3	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
4	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
5	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	90





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

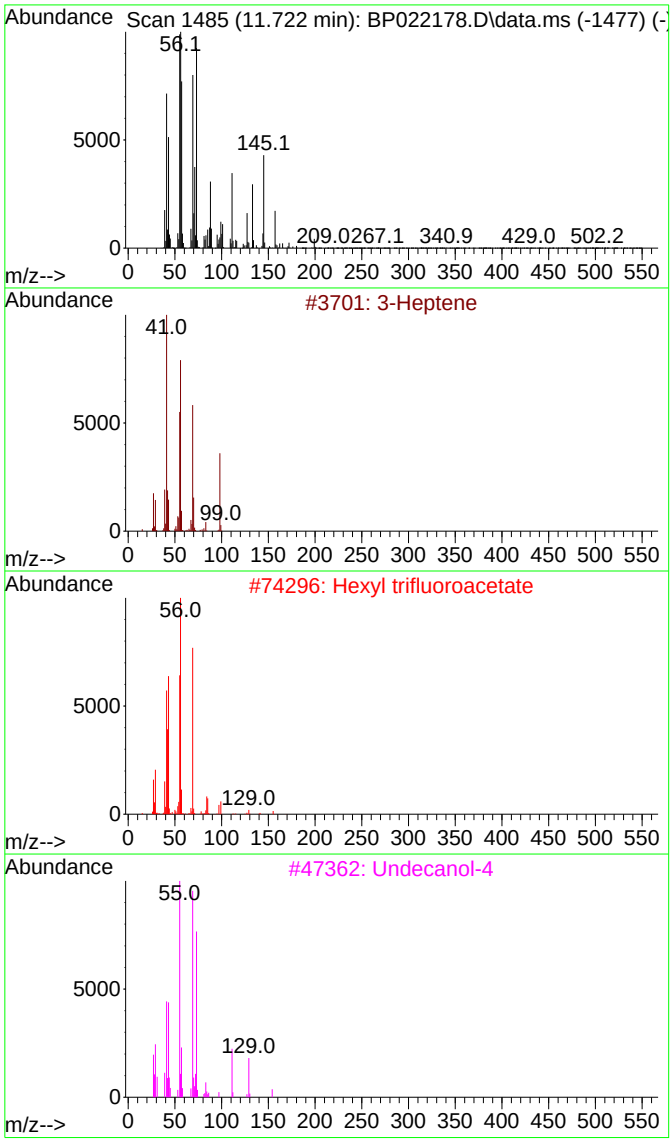
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

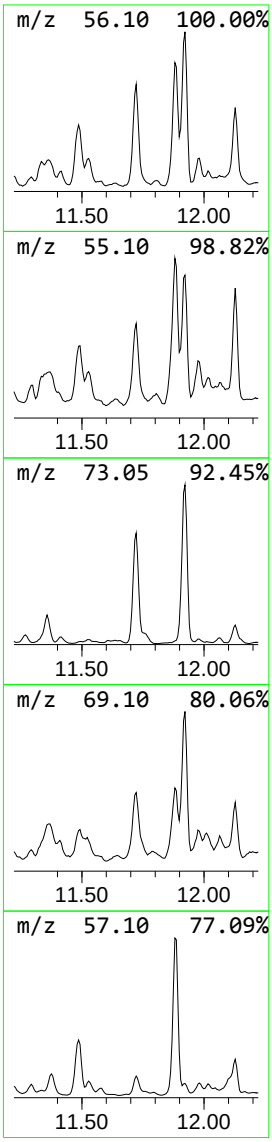
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.722	3.36 ng/ul	3000790	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptene	98	C7H14	000592-78-9	38
2	Hexyl trifluoroacetate	198	C8H13F3O2	000400-61-3	38
3	Undecanol-4	172	C11H24O	004272-06-4	22
4	1-Butanol, 3,3-dimethyl-	102	C6H14O	000624-95-3	22
5	Butyl caprylate	200	C12H24O2	000589-75-3	22





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

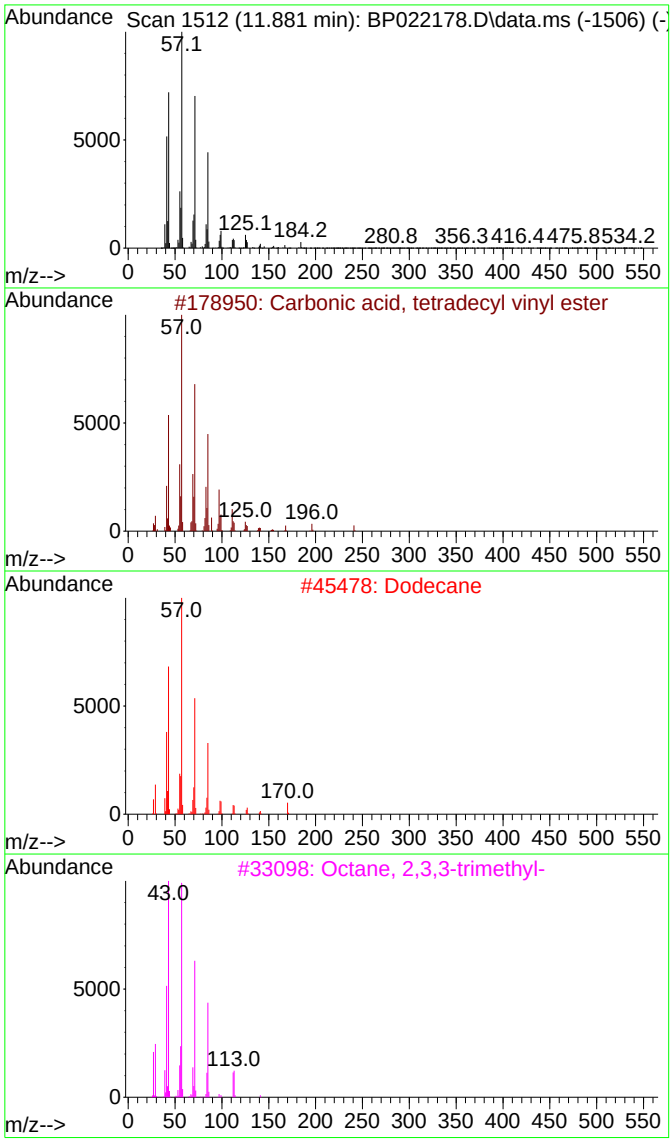
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

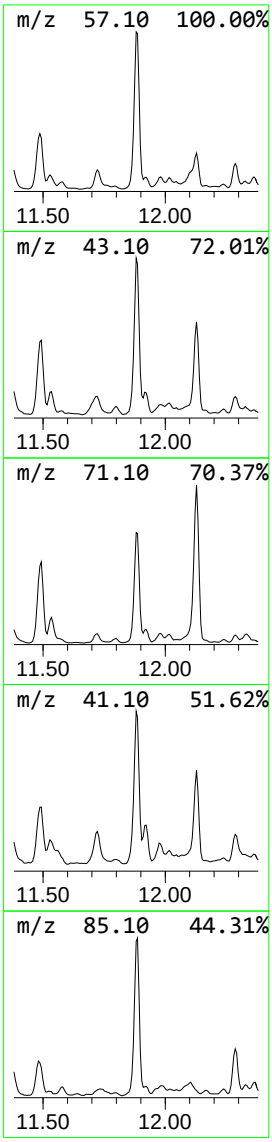
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 37 Carbonic acid, tetradecyl v... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.881	9.62 ng/ul	8582280	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	86
2	Dodecane	170	C12H26	000112-40-3	86
3	Octane, 2,3,3-trimethyl-	156	C11H24	062016-30-2	72
4	Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	72
5	Tridecane	184	C13H28	000629-50-5	70





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

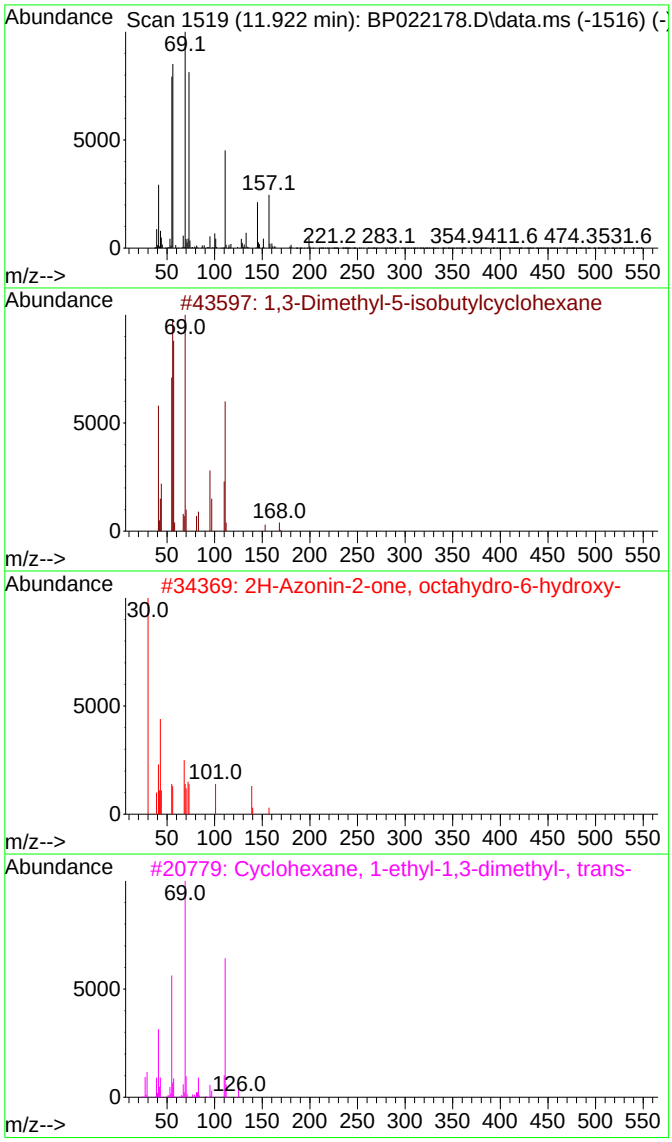
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

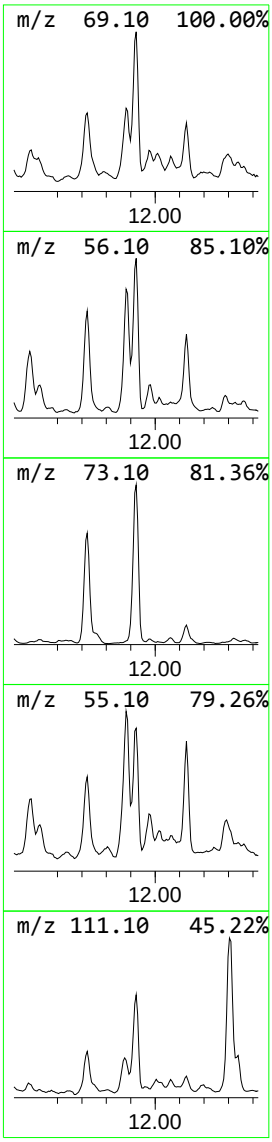
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.922	2.90 ng/ul	2583910	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dimethyl-5-isobutylcyclohexane	168	C12H24	013131-76-5	38
2	2H-Azonin-2-one, octahydro-6-hyd...	157	C8H15NO2	023435-07-6	38
3	Cyclohexane, 1-ethyl-1,3-dimethy...	140	C10H20	062238-29-3	17
4	Cyclohexane, 1-ethyl-1,4-dimethy...	140	C10H20	062238-32-8	17
5	Cyclohexane, 1-ethyl-1,3-dimethy...	140	C10H20	062238-31-7	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

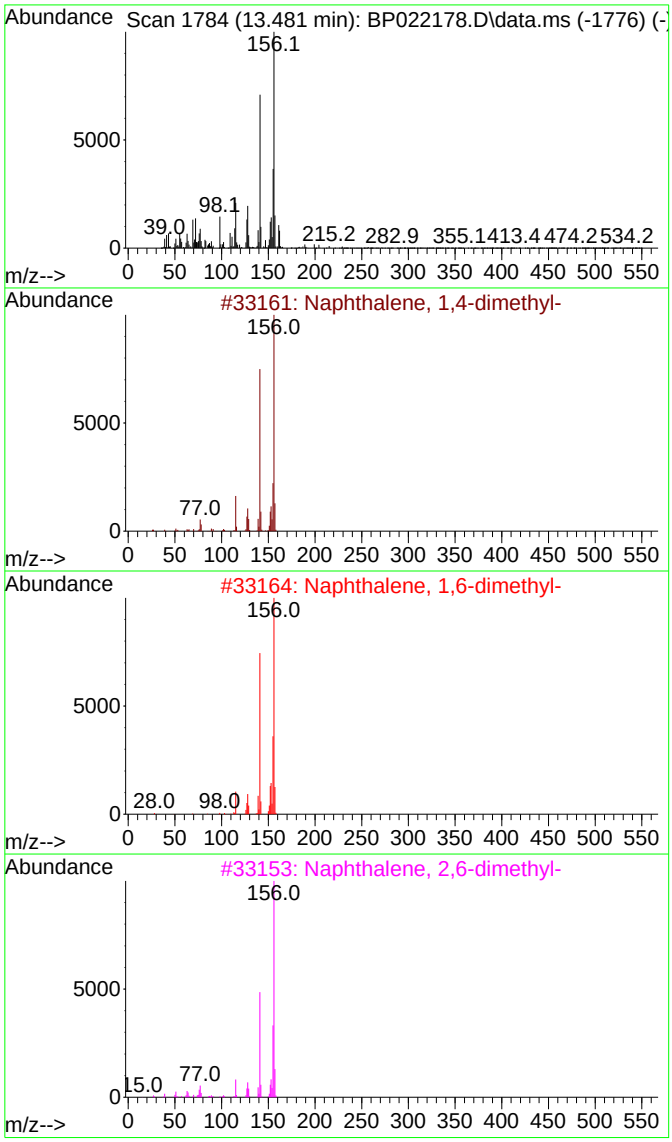
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

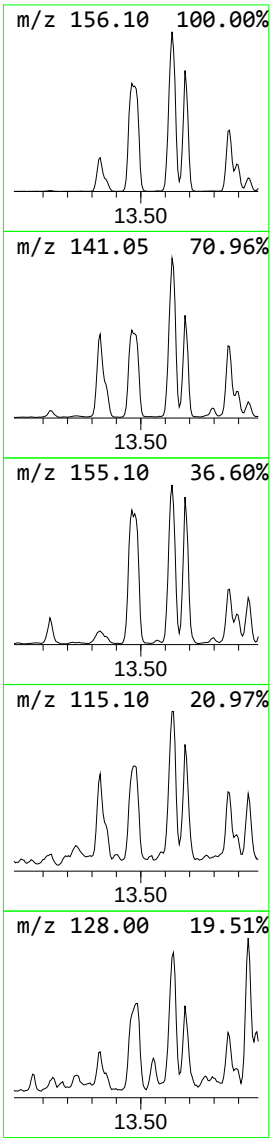
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 41 Naphthalene, 1,4-dimethyl- Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.481	3.98 ng/ul	6540660	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

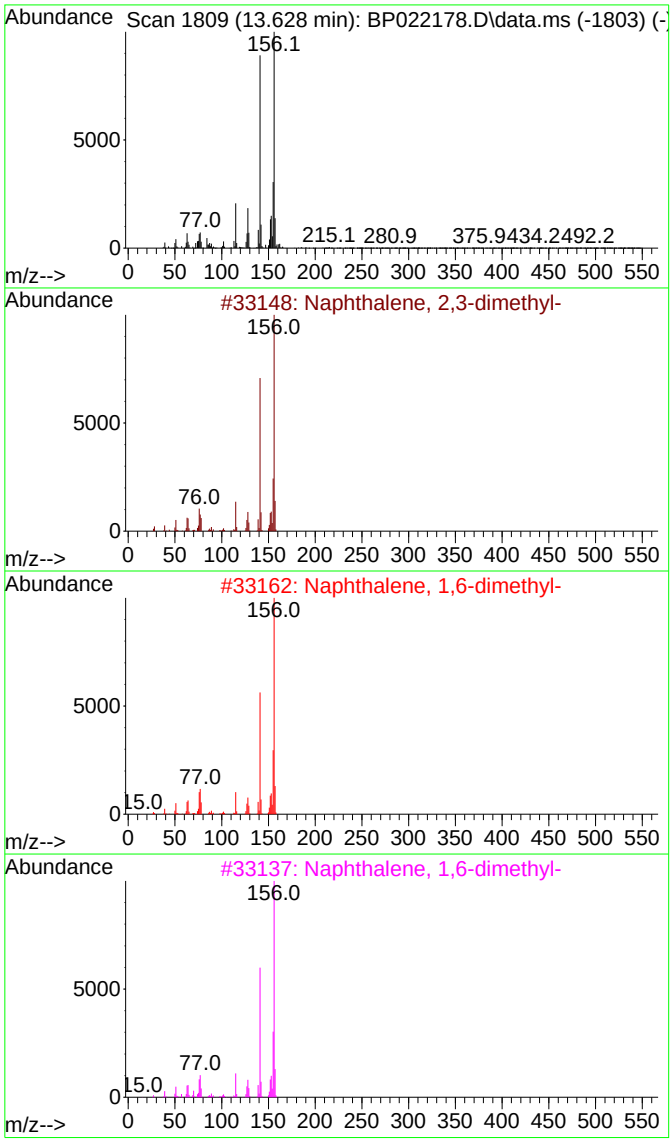
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

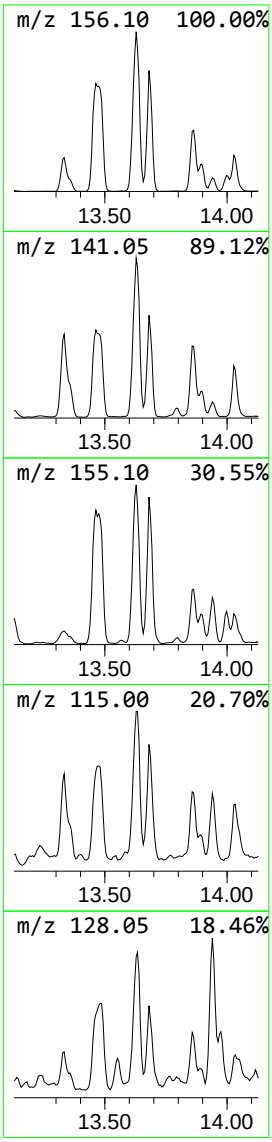
****

Peak Number 42 Naphthalene, 2,3-dimethyl- Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.628	3.93 ng/ul	6454700	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

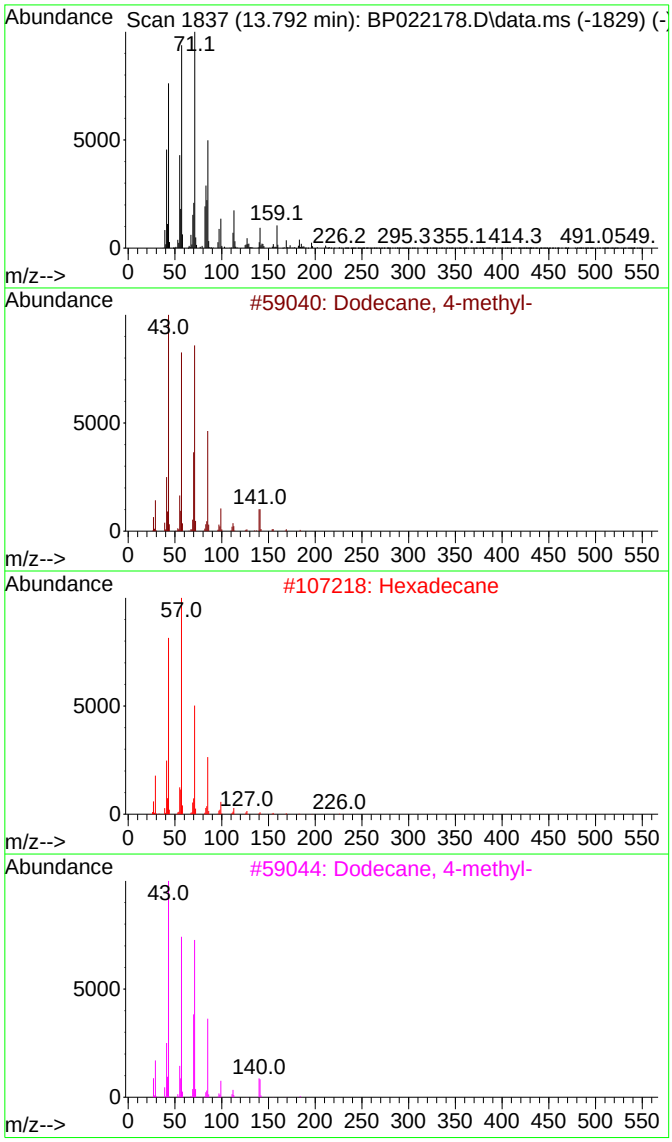
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

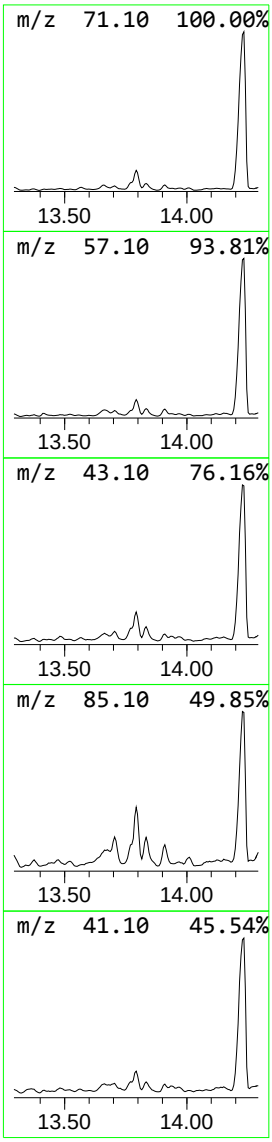
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.792	2.85 ng/ul	4684170	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 4-methyl-	184	C13H28	006117-97-1	60
2	Hexadecane	226	C16H34	000544-76-3	53
3	Dodecane, 4-methyl-	184	C13H28	006117-97-1	53
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	52
5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	50





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

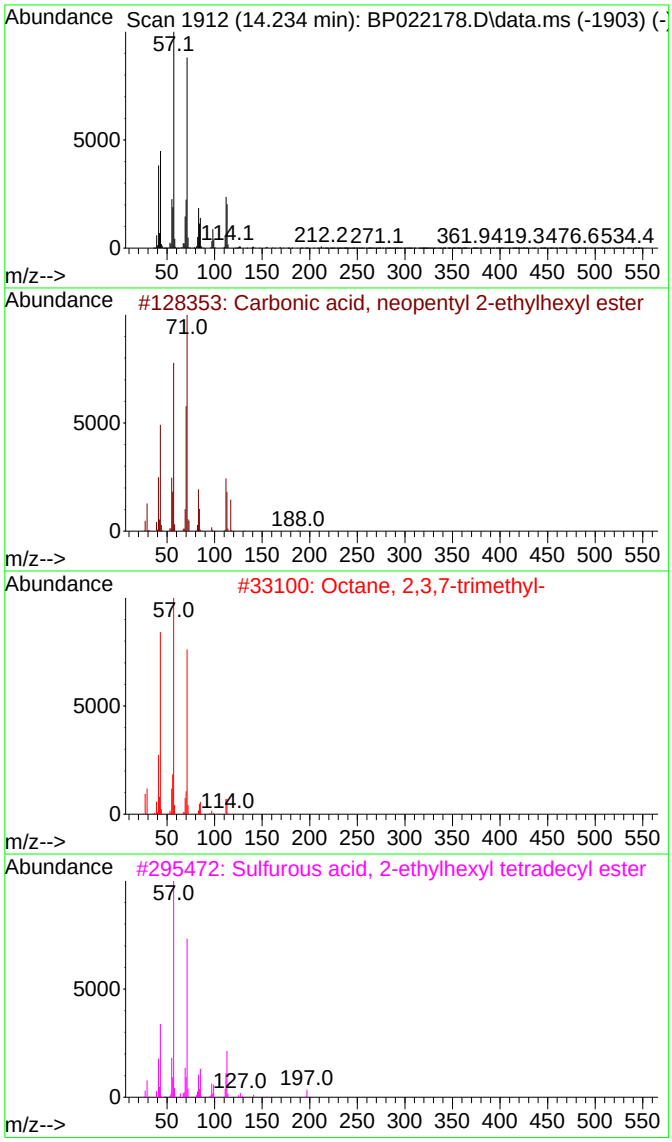
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

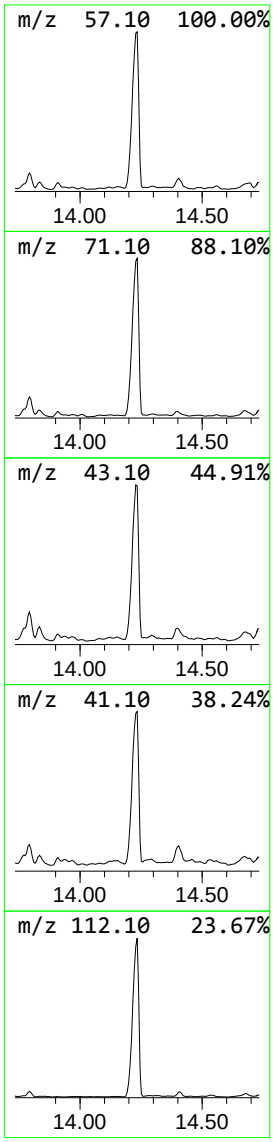
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 45 Carbonic acid, neopentyl 2-... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.234	20.00 ng/ul	32862600	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	74
2	Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
3	Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	64
4	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
5	Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	59





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

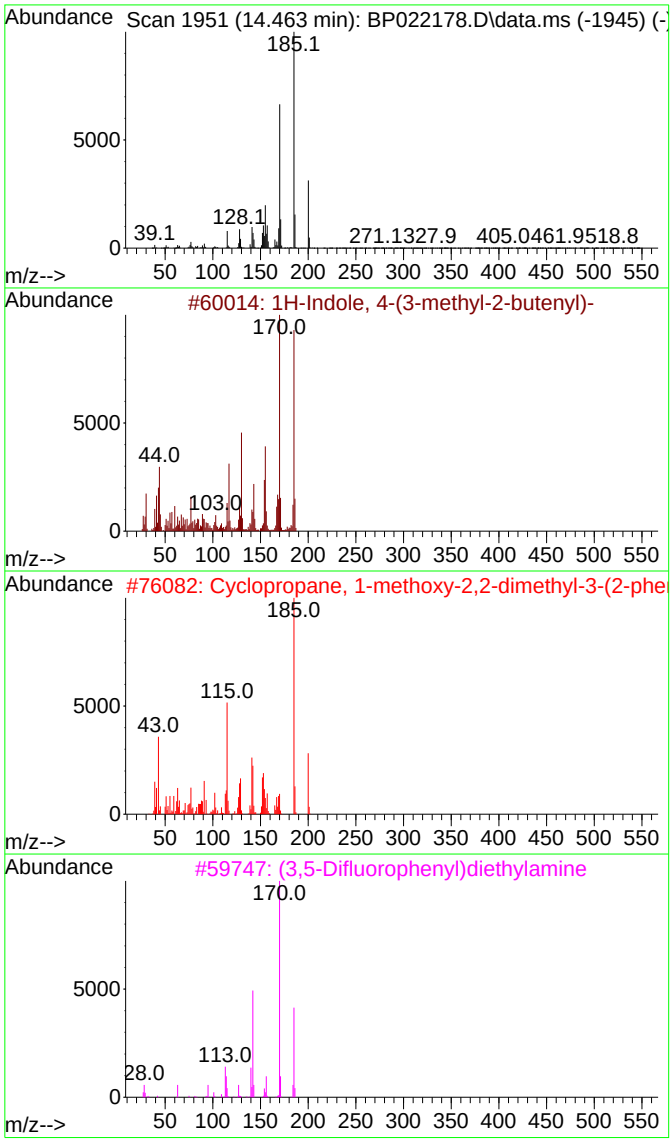
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

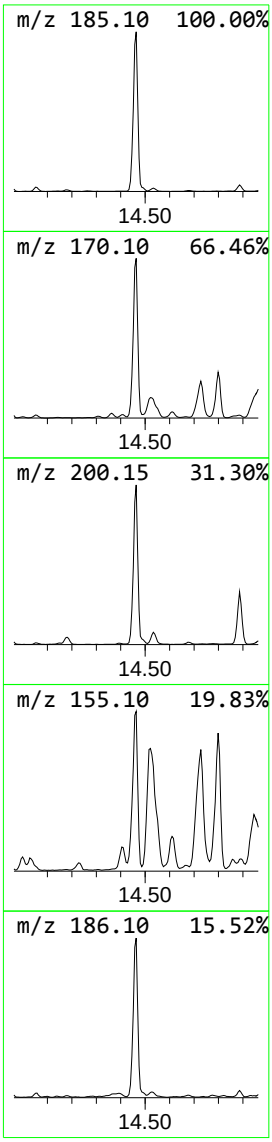
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 47 1H-Indole, 4-(3-methyl-2-bu... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	8.70 ng/ul	14291700	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indole, 4-(3-methyl-2-butenyl)-	185	C13H15N	032962-25-7	72
2	Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	58
3	(3,5-Difluorophenyl)diethylamine	185	C10H13F2N	1000306-62-9	52
4	Ethanone, 1-(6-methoxy-1-naphtha...	200	C13H12O2	058149-89-6	47
5	1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	46





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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

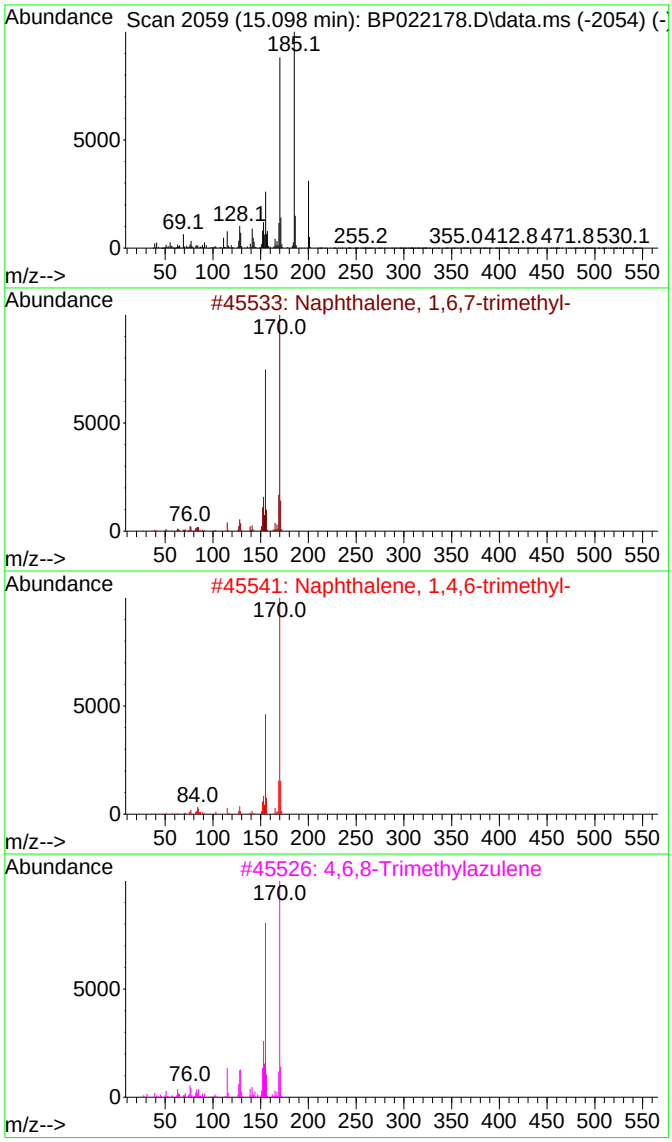
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

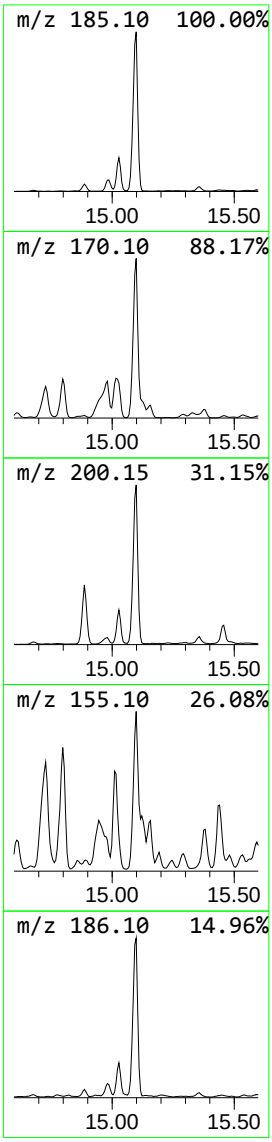
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 49 Naphthalene, 1,6,7-trimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.098	10.00 ng/ul	16427600	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	50
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
3	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	46
4	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	46
5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	45





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Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

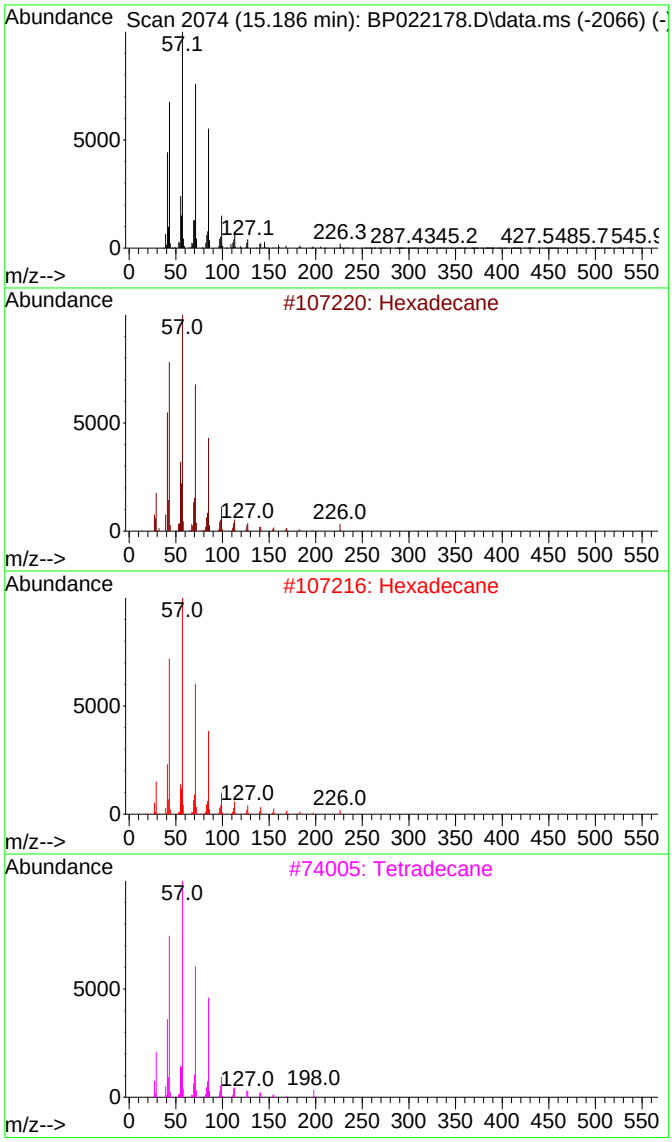
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

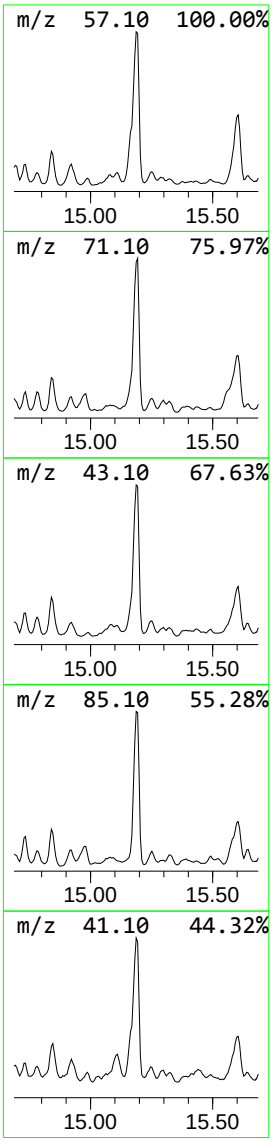
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.187	7.77 ng/ul	12772200	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	95
2	Hexadecane	226	C16H34	000544-76-3	94
3	Tetradecane	198	C14H30	000629-59-4	91
4	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

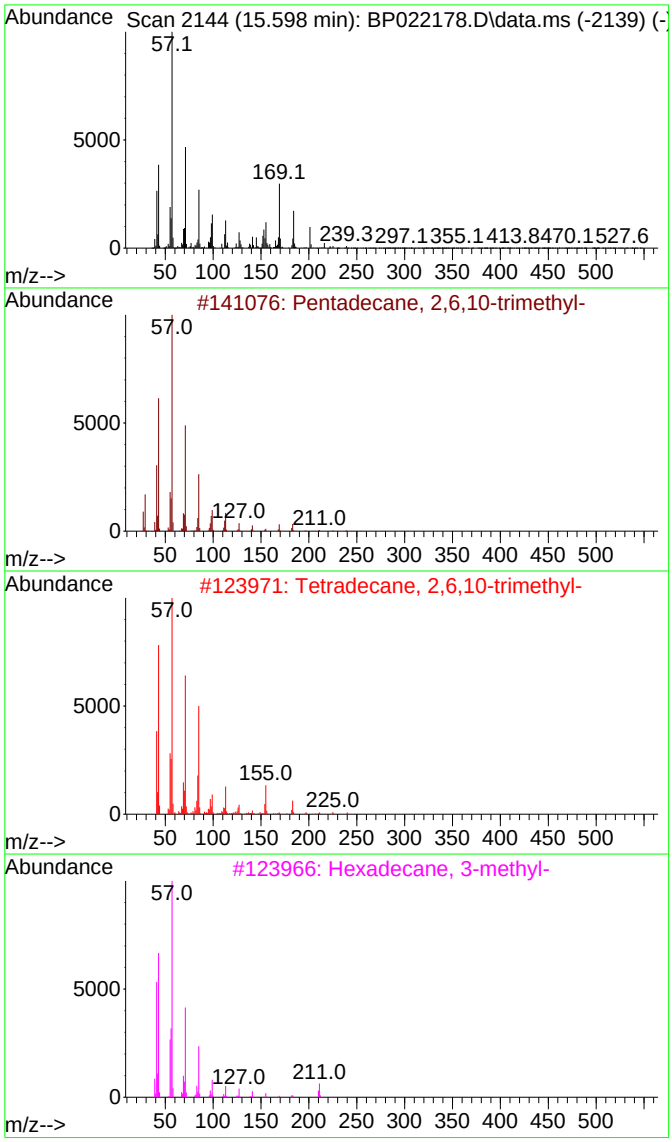
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

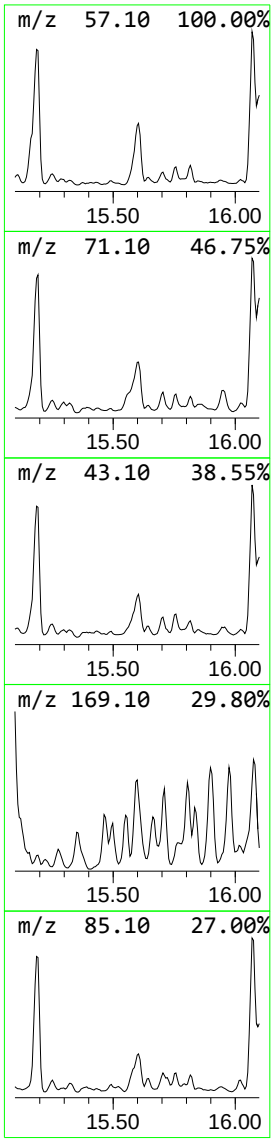
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 64

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.598	2.71 ng/ul	4454330	Acenaphthene-d10	14.310

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	78
2	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	49
3	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	49
4	Hexadecane	226	C16H34	000544-76-3	46
5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	43





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

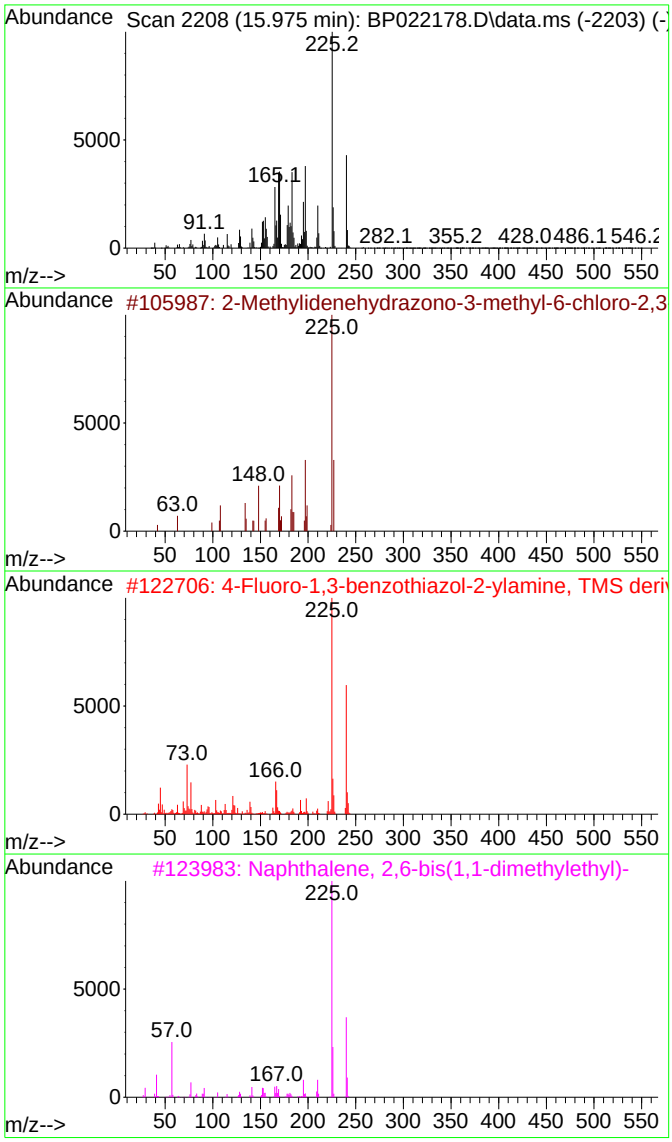
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

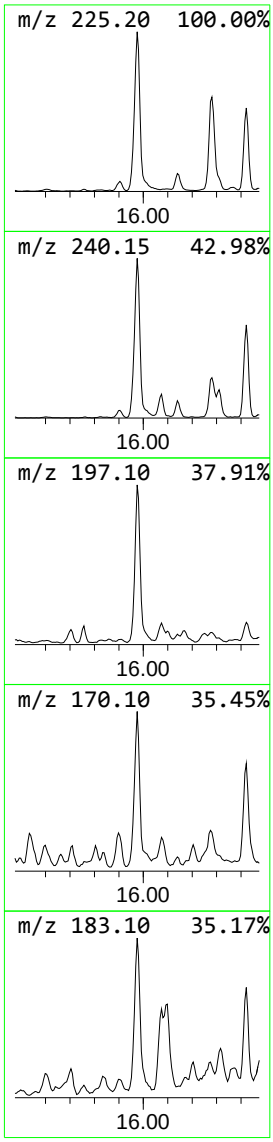
****

Peak Number 53 unknown-04 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.975	13.99 ng/ul	4058370	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylidenehydrazono-3-methyl-...	225	C9H8ClN3S	1000147-91-5	43
2	4-Fluoro-1,3-benzothiazol-2-ylam...	240	C10H13FN2SSi	1000478-75-9	38
3	Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	38
4	2-Methylidene-hydrazono-3-methyl...	225	C9H8ClN3S	1000147-91-3	35
5	[1-(1,5,5-Trimethyl-2,4-dioxopyr...	240	C11H16N2O4	1010287-56-8	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

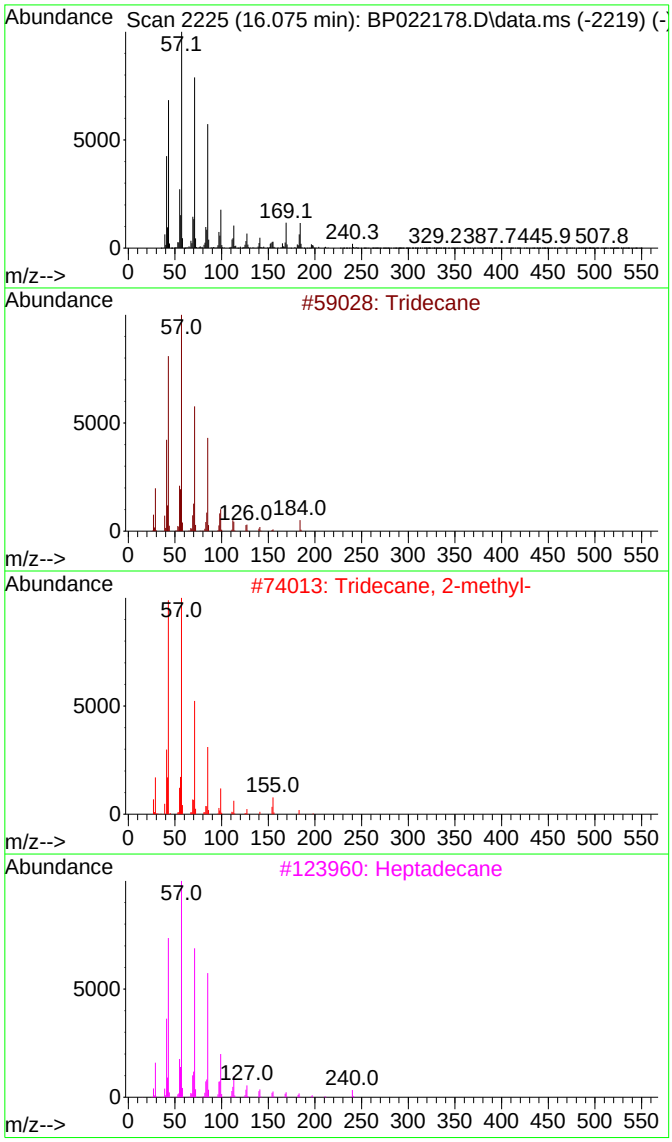
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

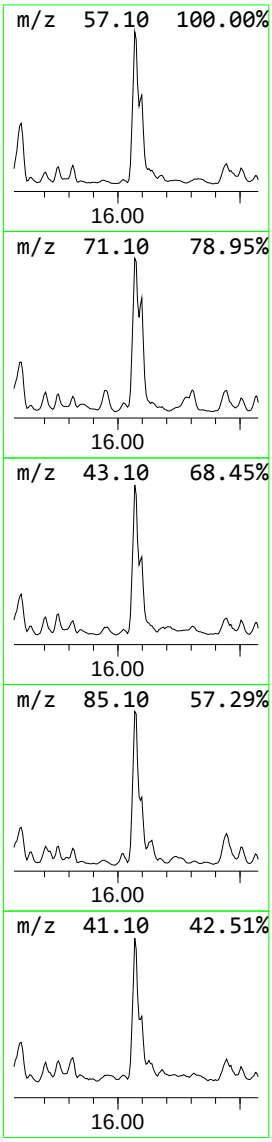
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.075	46.75 ng/ul	13561100	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	90
2	Tridecane, 2-methyl-	198	C14H30	001560-96-9	87
3	Heptadecane	240	C17H36	000629-78-7	86
4	Octadecane	254	C18H38	000593-45-3	83
5	Heneicosane	296	C21H44	000629-94-7	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

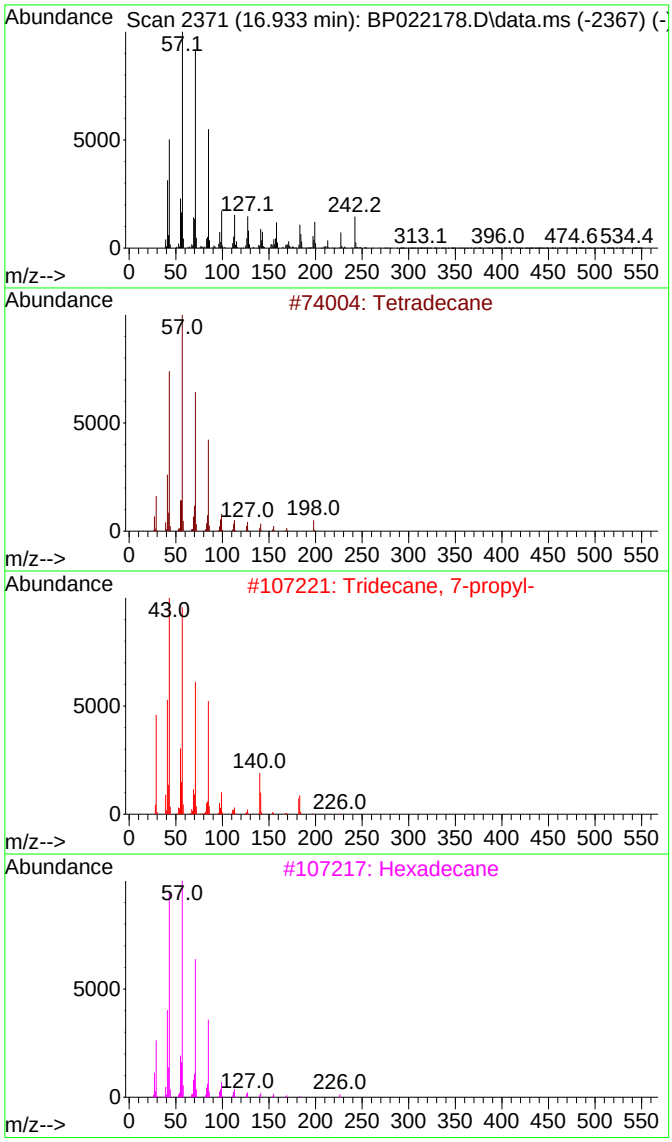
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

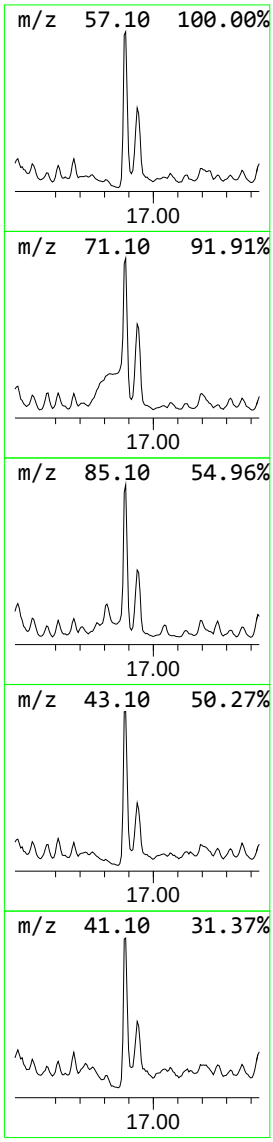
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.933	28.72 ng/ul	8330520	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	62
2	Tridecane, 7-propyl-	226	C16H34	055045-09-5	60
3	Hexadecane	226	C16H34	000544-76-3	58
4	Dodecane	170	C12H26	000112-40-3	58
5	2-Bromotetradecane	276	C14H29Br	074036-95-6	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

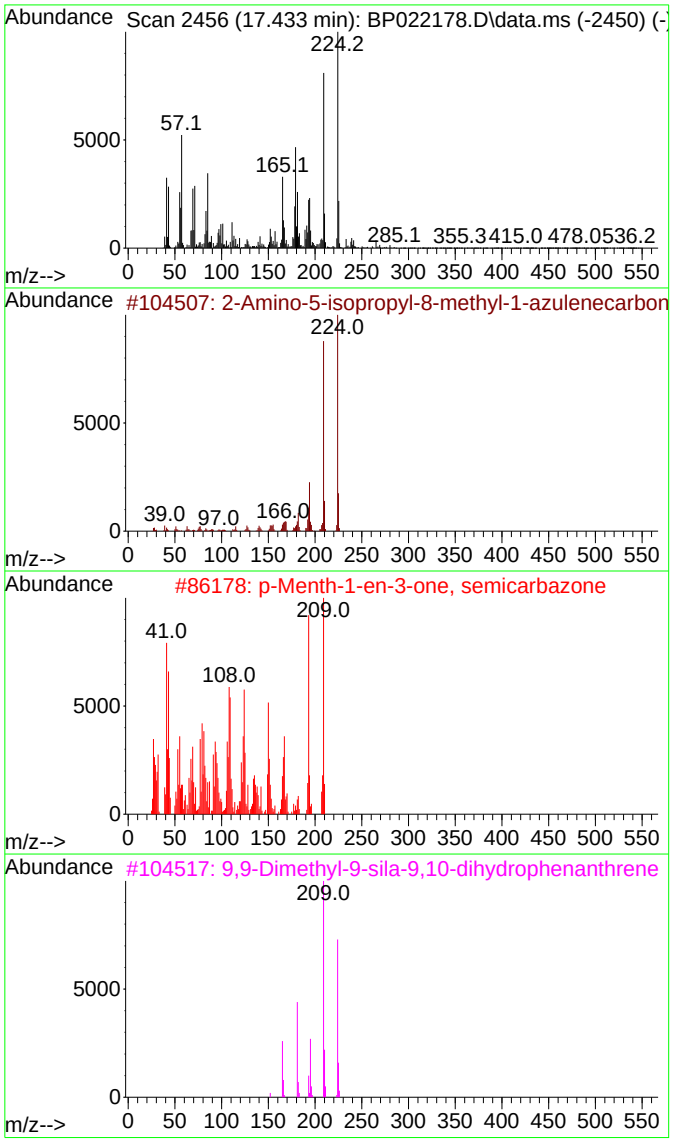
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

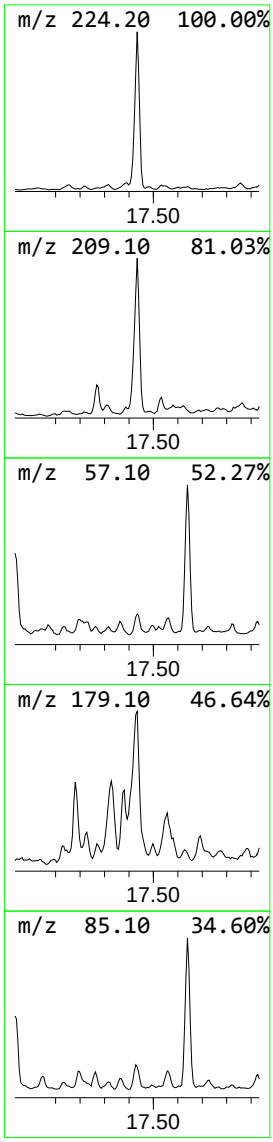
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 56 2-Amino-5-isopropyl-8-methy... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.433	19.05 ng/ul	5526090	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-5-isopropyl-8-methyl-1-a...	224	C15H16N2	093946-48-6	83
2	p-Menth-1-en-3-one, semicarbazone	209	C11H19N3O	004713-41-1	74
3	9,9-Dimethyl-9-sila-9,10-dihydro...	224	C15H16Si	187328-55-8	58
4	1H-1,3-Benzimidazole, 5-methoxy-...	224	C14H12N2O	1000350-62-4	53
5	Methyl 4,6-bis(dimethylamino)-1,...	224	C9H16N6O	1000101-98-0	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

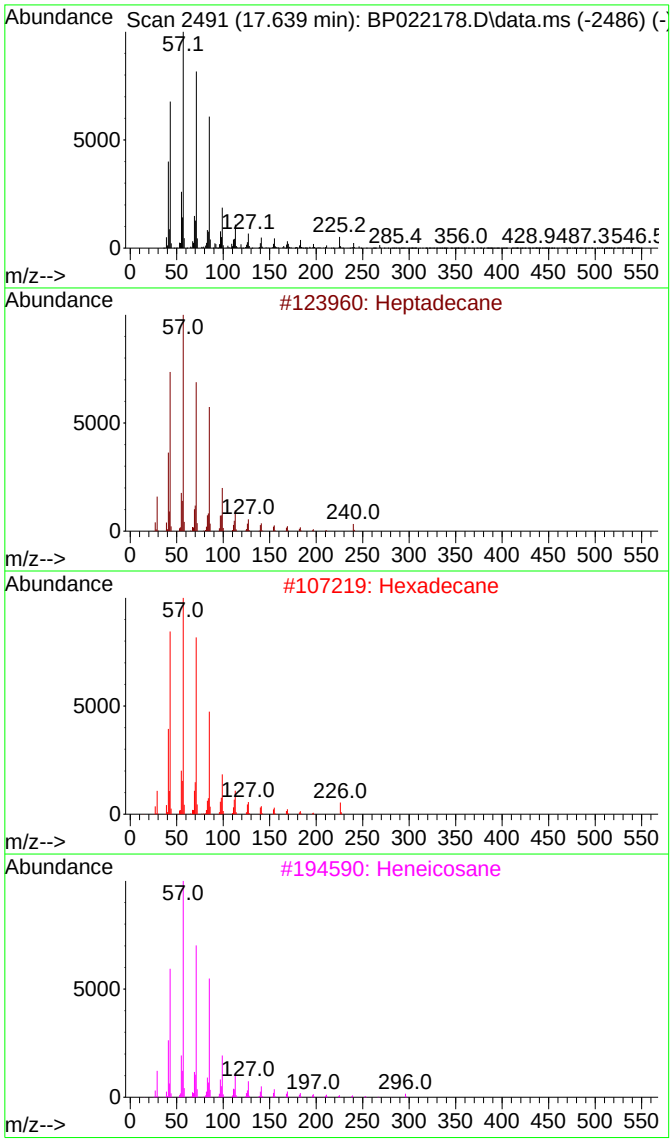
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

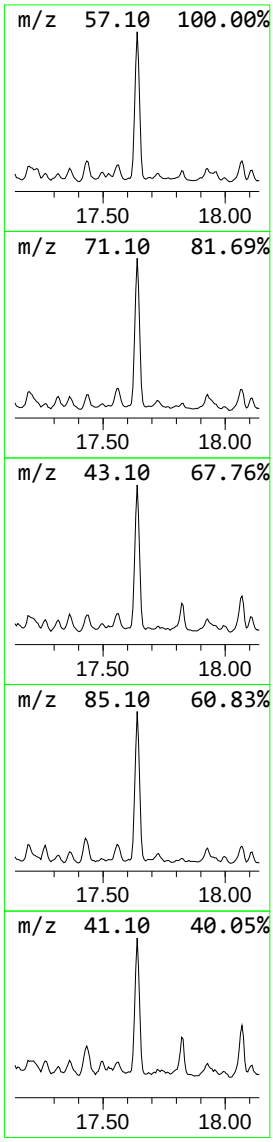
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.639	30.05 ng/ul	8716900	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	97
2	Hexadecane	226	C16H34	000544-76-3	93
3	Heneicosane	296	C21H44	000629-94-7	90
4	Octadecane	254	C18H38	000593-45-3	90
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

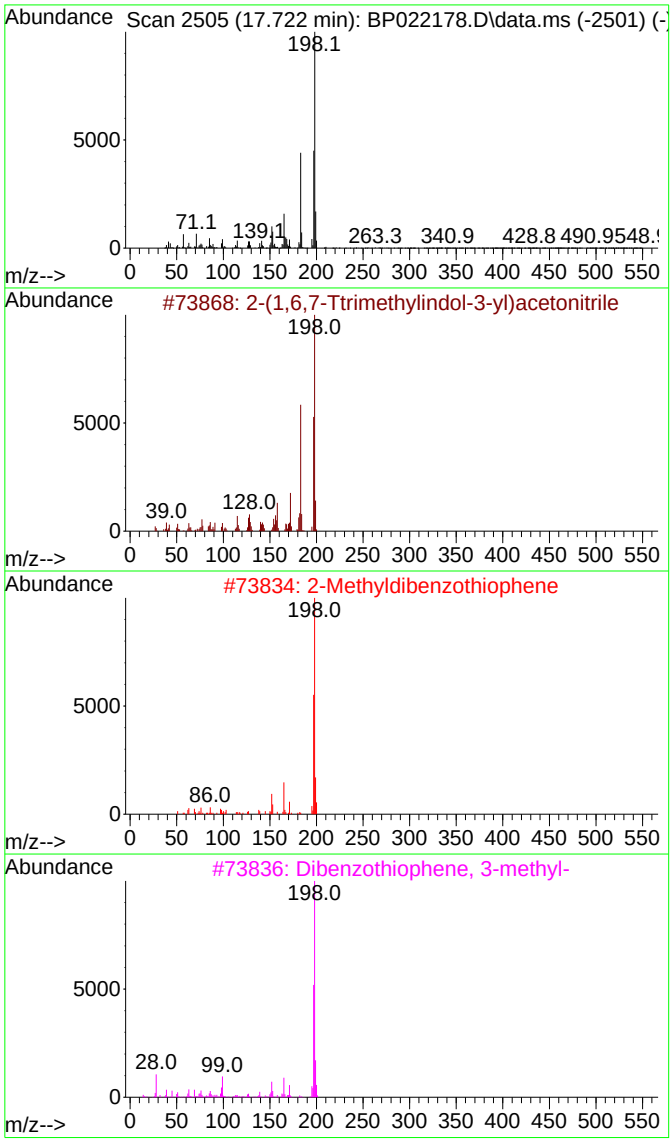
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

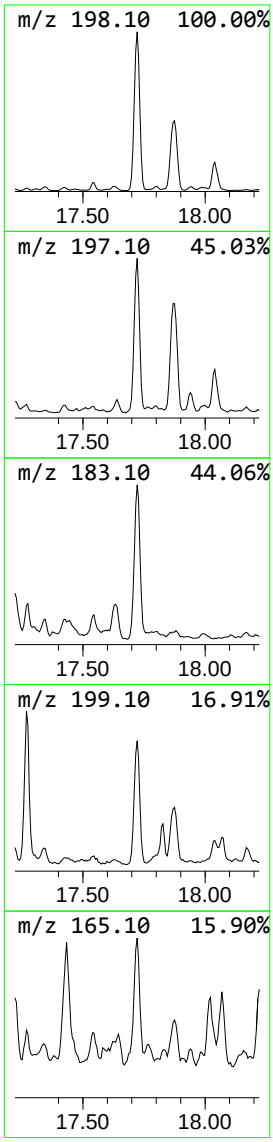
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 58 2-(1,6,7-Ttrimethylindol-3-... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.722	9.62 ng/ul	2790250	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(1,6,7-Ttrimethylindol-3-yl)ac...	198	C13H14N2	1000436-09-3	80
2	2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	76
3	Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	68
4	Benzene, 1-methyl-2-(4-methylphe...	198	C14H14O	003402-72-0	59
5	4,6,8-Trimethyl-1-azulenecarbald...	198	C14H14O	000832-62-2	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
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Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

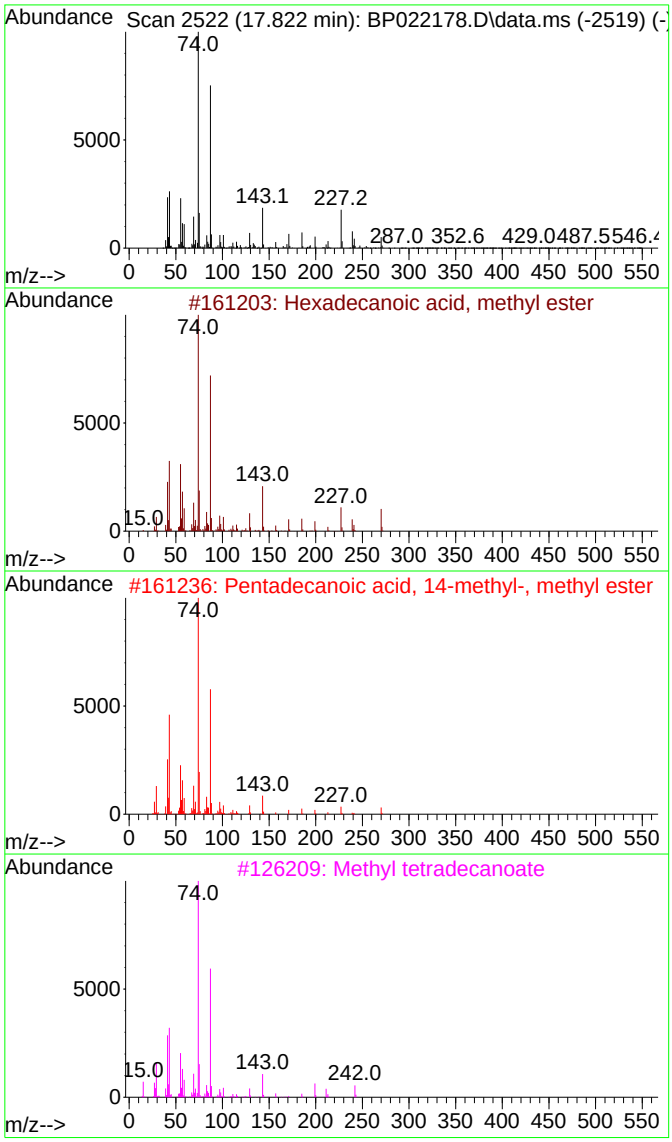
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

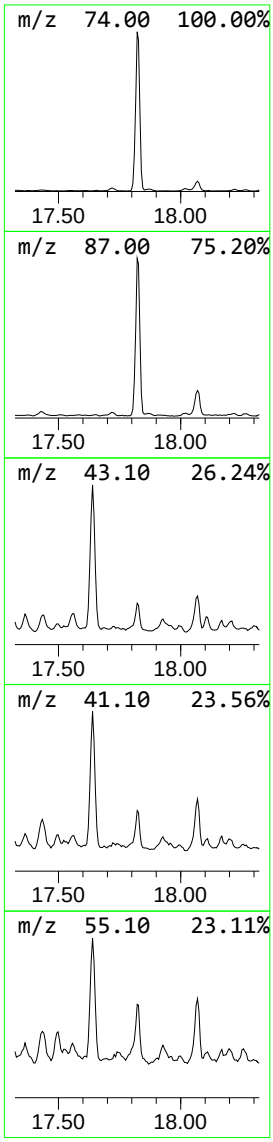
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 59 Hexadecanoic acid, methyl e... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.822	8.76 ng/ul	2539940	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
2	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3	Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
4	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	87





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

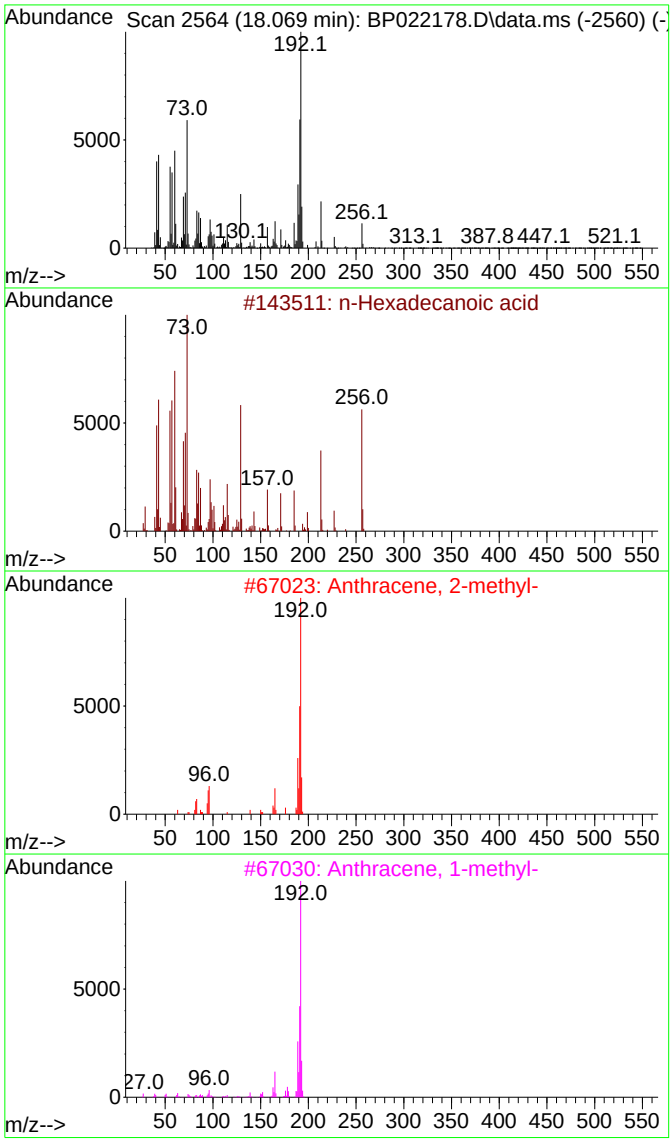
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

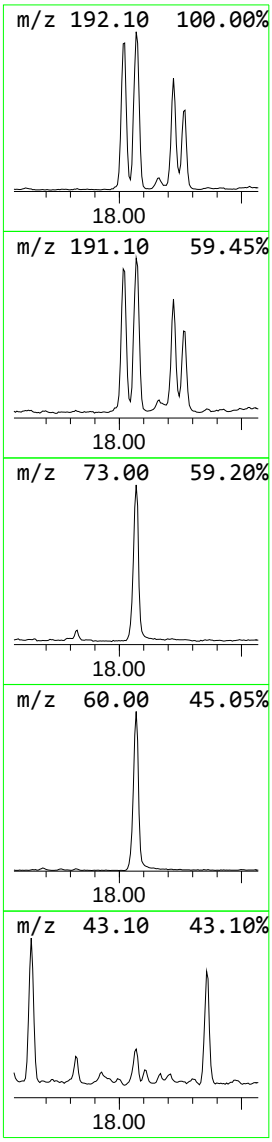
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 60 n-Hexadecanoic acid Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.069	14.88 ng/ul	4316560	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	70
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	70





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

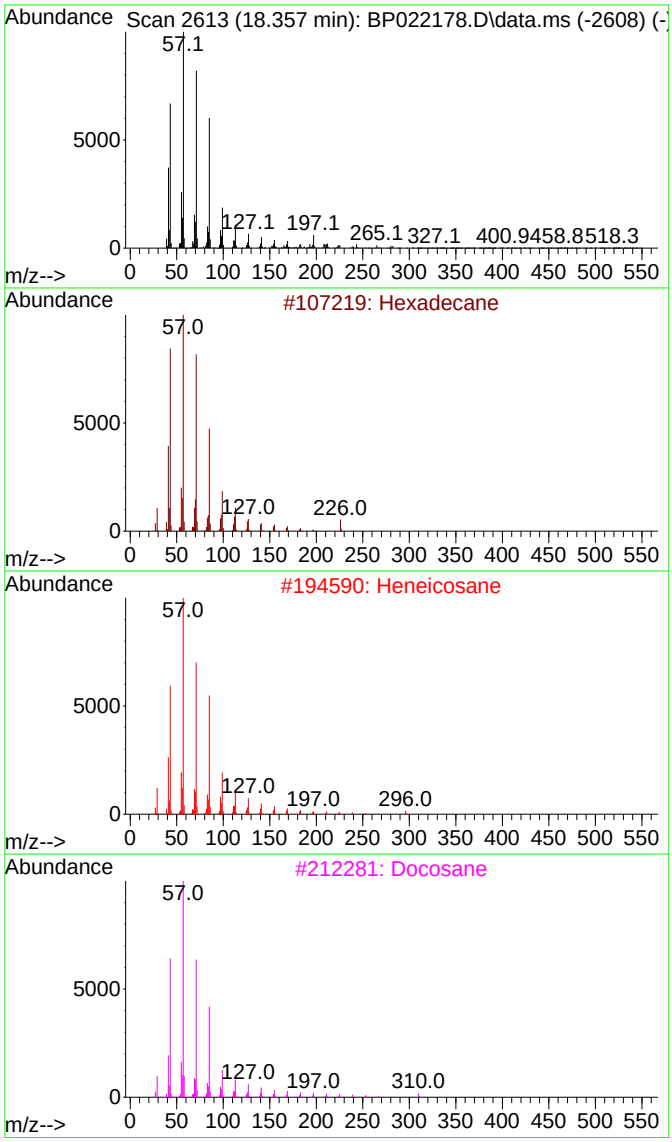
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

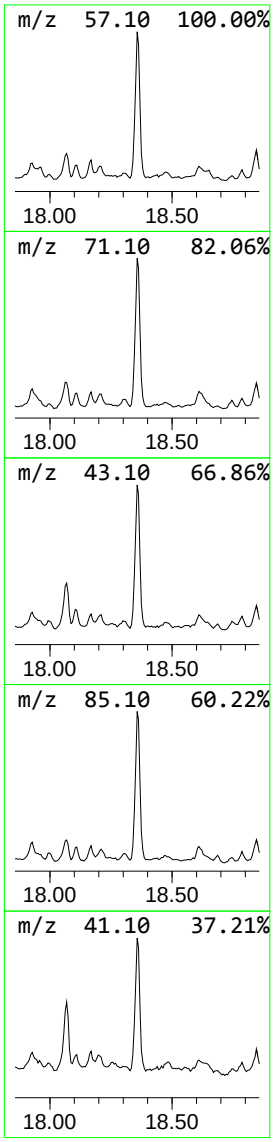
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.357	23.73 ng/ul	6882260	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Heneicosane	296	C21H44	000629-94-7	90
3	Docosane	310	C22H46	000629-97-0	87
4	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
5	Octadecane	254	C18H38	000593-45-3	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

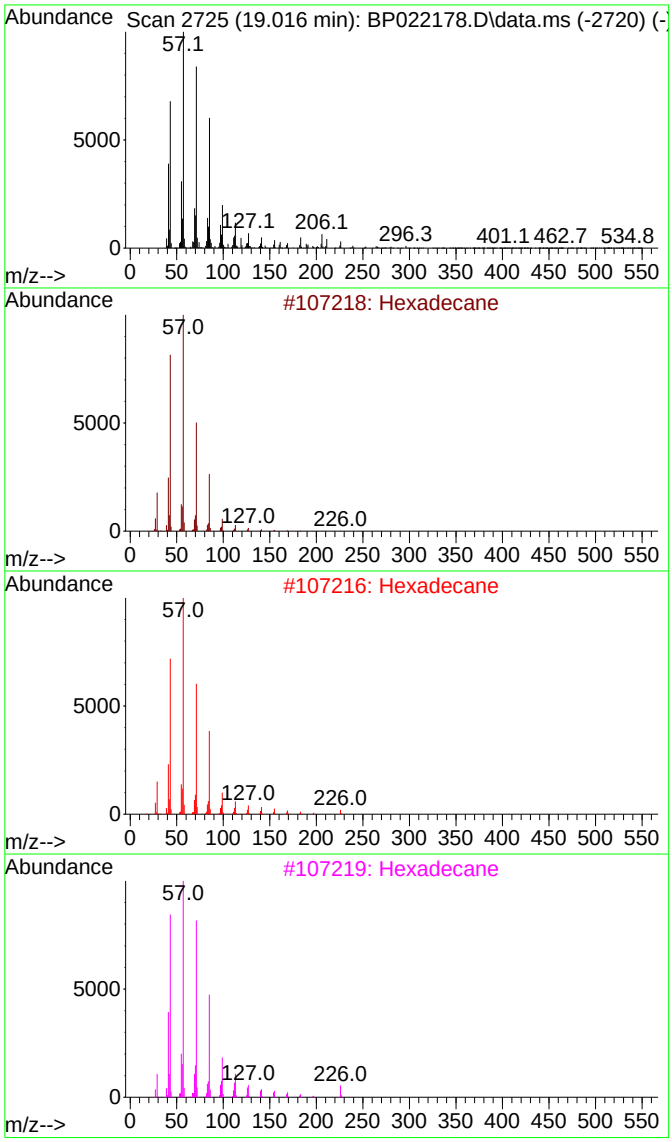
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

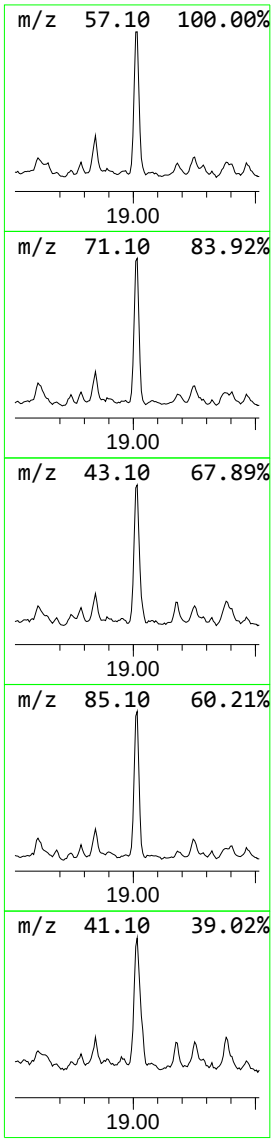
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.016	27.18 ng/ul	7882560	Phenanthrene-d10	17.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Hexadecane	226	C16H34	000544-76-3	93
3	Hexadecane	226	C16H34	000544-76-3	93
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
5	Heneicosane	296	C21H44	000629-94-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

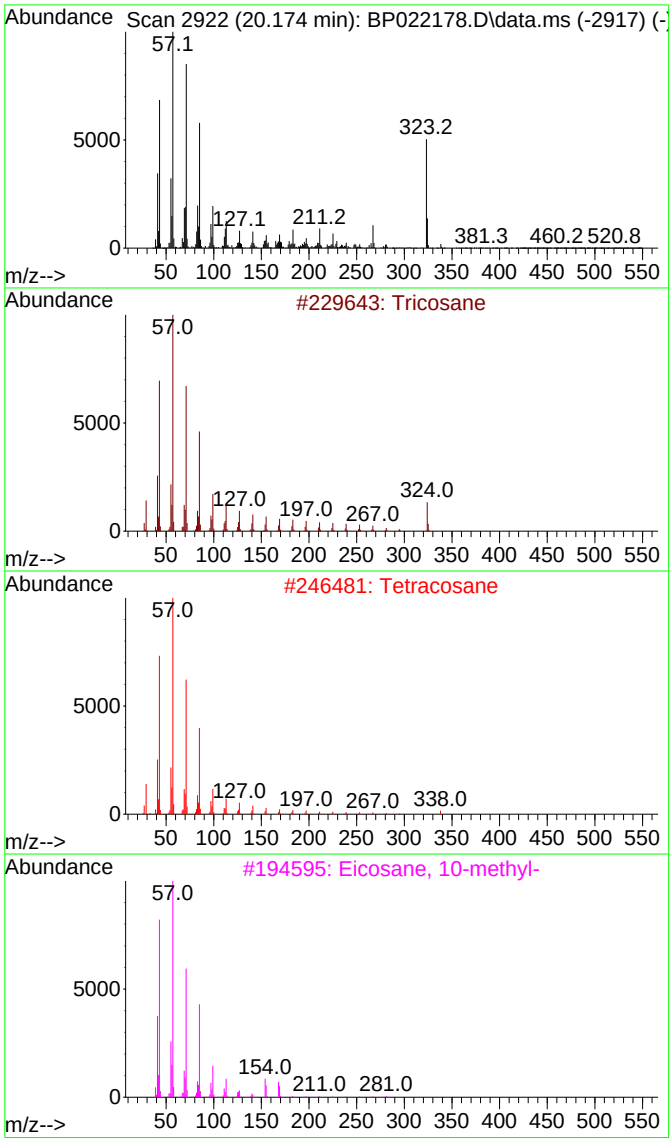
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

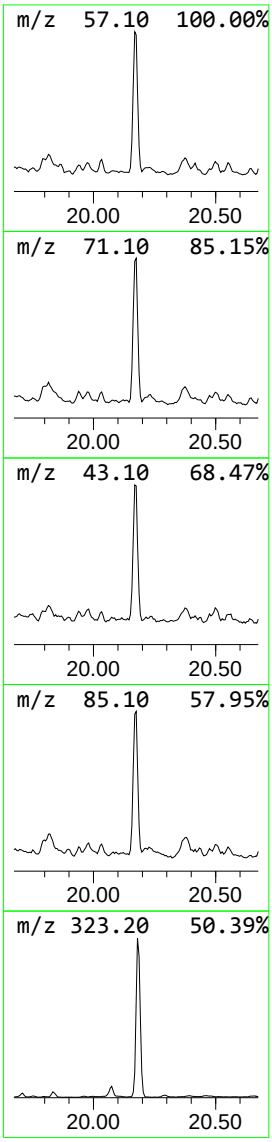
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.174	33.97 ng/ul	4696110	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	92
2	Tetracosane	338	C24H50	000646-31-1	87
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	83
4	Tetracosane	338	C24H50	000646-31-1	83
5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

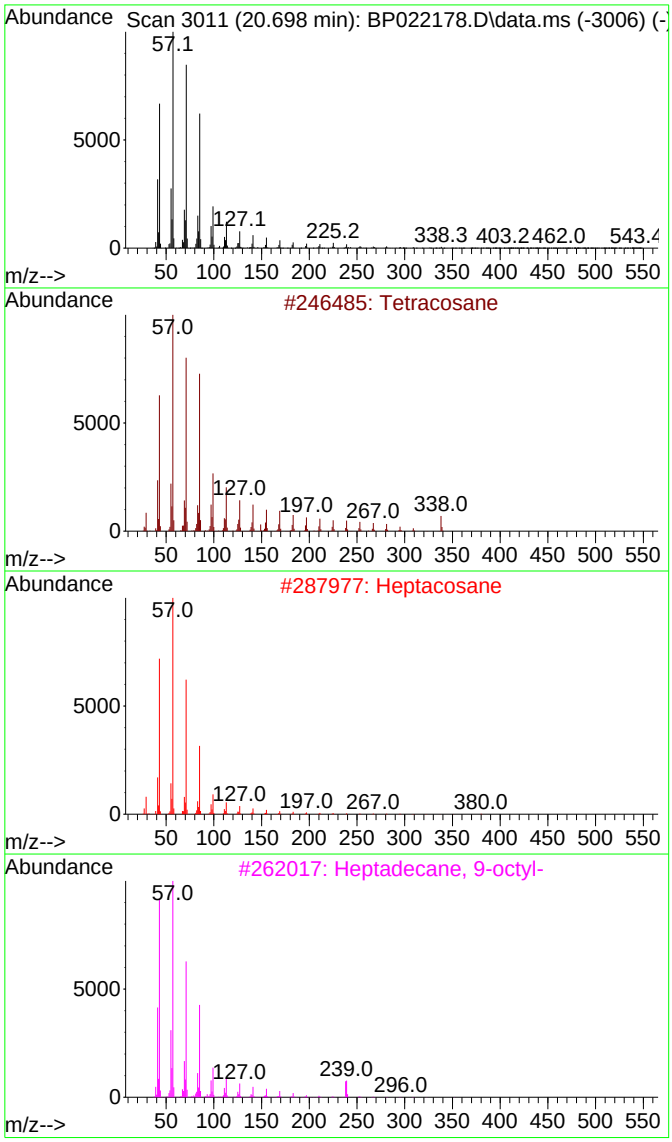
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

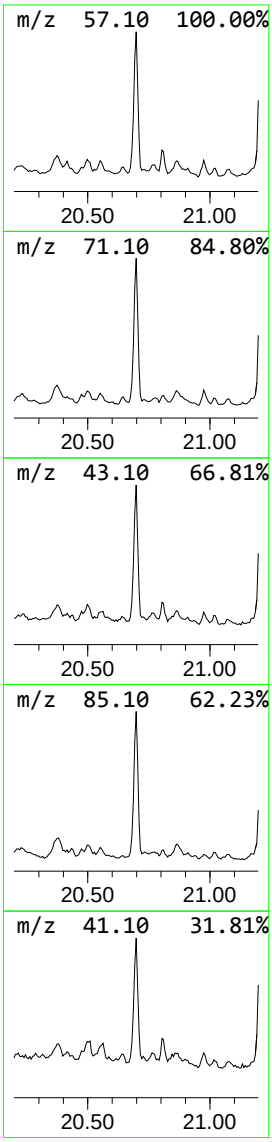
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.698	19.73 ng/ul	2727850	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	94
2	Heptacosane	380	C27H56	000593-49-7	91
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Docosane	310	C22H46	000629-97-0	91
5	Heneicosane	296	C21H44	000629-94-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

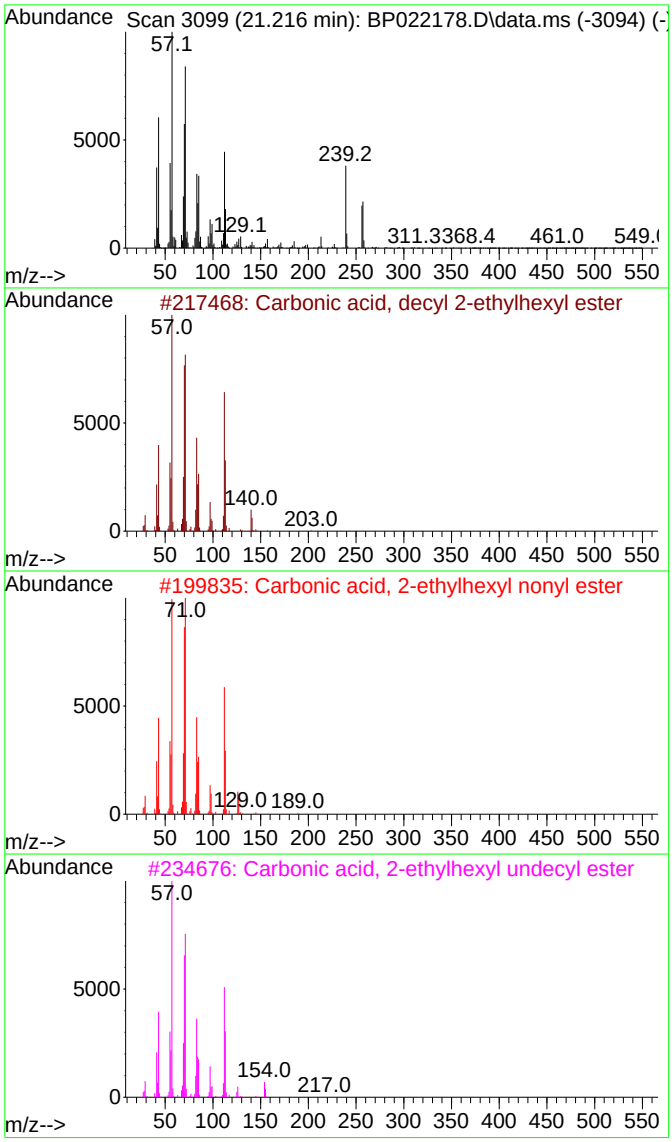
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

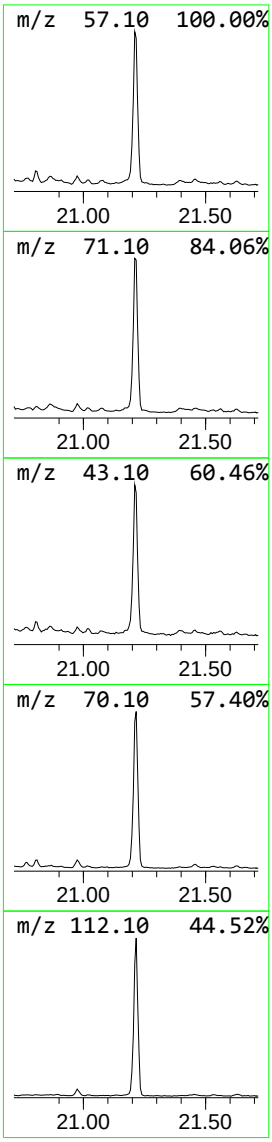
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 65 Carbonic acid, decyl 2-ethy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.216	64.90 ng/ul	8971370	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, decyl 2-ethylhexy...	314	C19H38O3	1000383-13-7	58
2	Carbonic acid, 2-ethylhexyl nony...	300	C18H36O3	1000383-13-6	58
3	Carbonic acid, 2-ethylhexyl unde...	328	C20H40O3	1000383-13-8	52
4	CITRONELLAL, DIHYDRO-	156	C10H20O	005988-91-0	46
5	Carbonic acid, 2-ethylhexyl octy...	286	C17H34O3	1000383-13-5	43





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Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

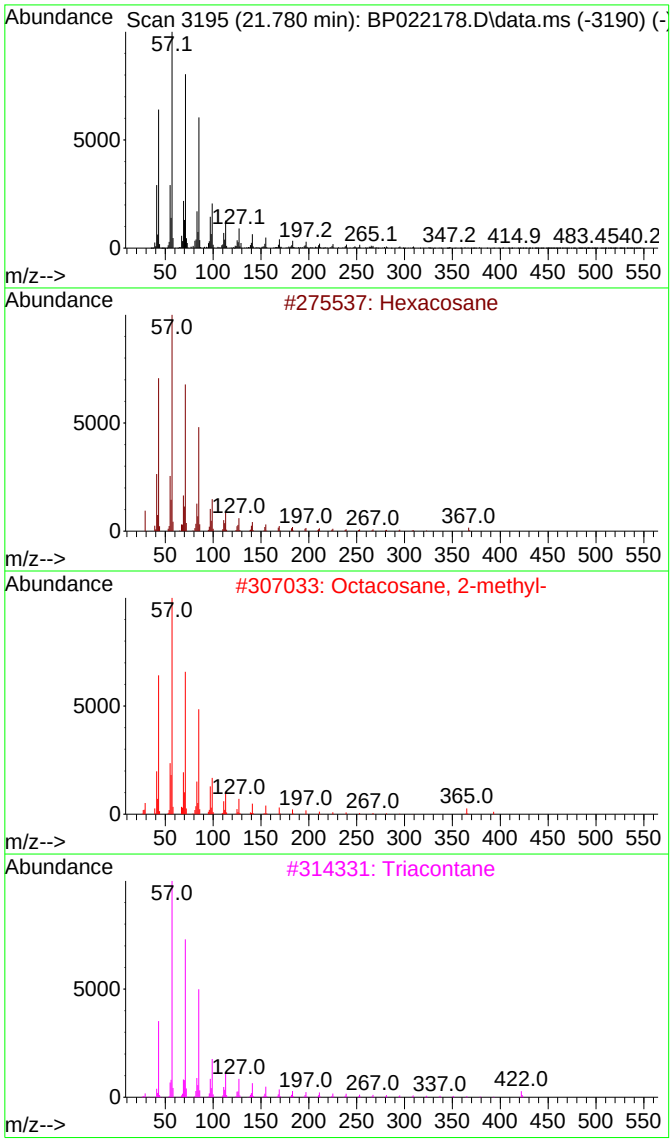
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

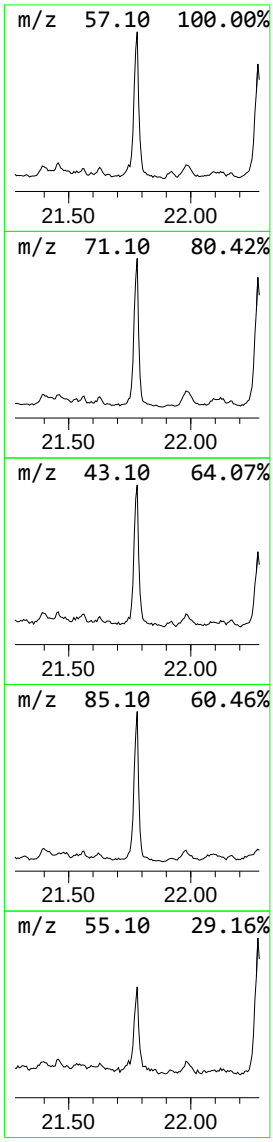
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.780	25.41 ng/ul	3512360	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	94
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3	Triacontane	422	C30H62	000638-68-6	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Tetracosane	338	C24H50	000646-31-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

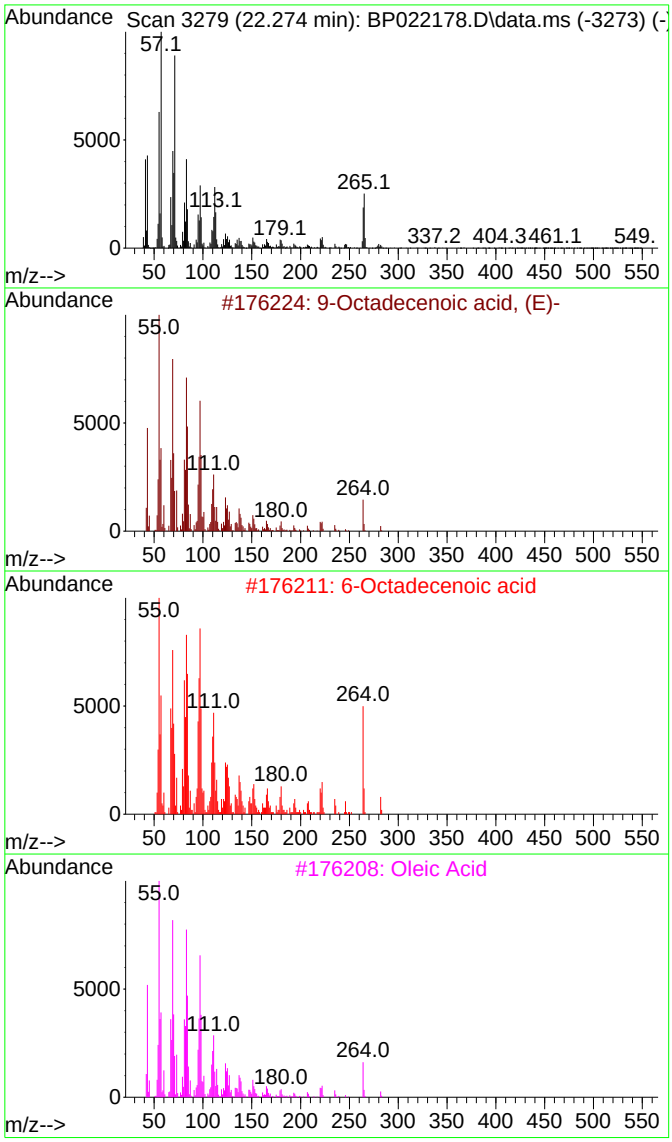
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

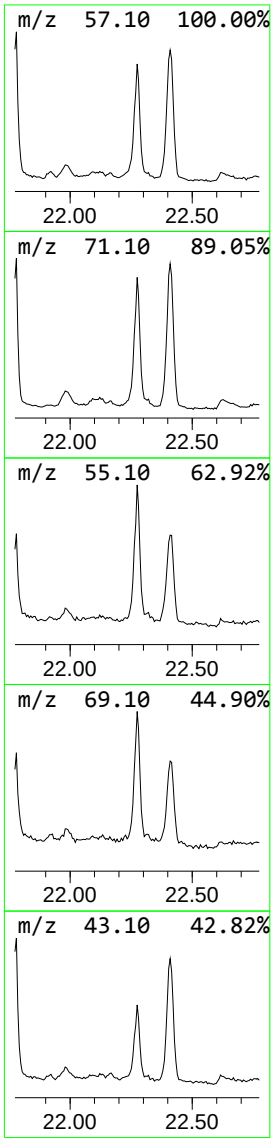
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 67 9-Octadecenoic acid, (E)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.274	29.18 ng/ul	4034210	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	56
2	6-Octadecenoic acid	282	C18H34O2	1000336-66-8	53
3	Oleic Acid	282	C18H34O2	000112-80-1	44
4	Oxalic acid, 2-ethylhexyl pentad...	412	C25H48O4	1000309-39-8	43
5	Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

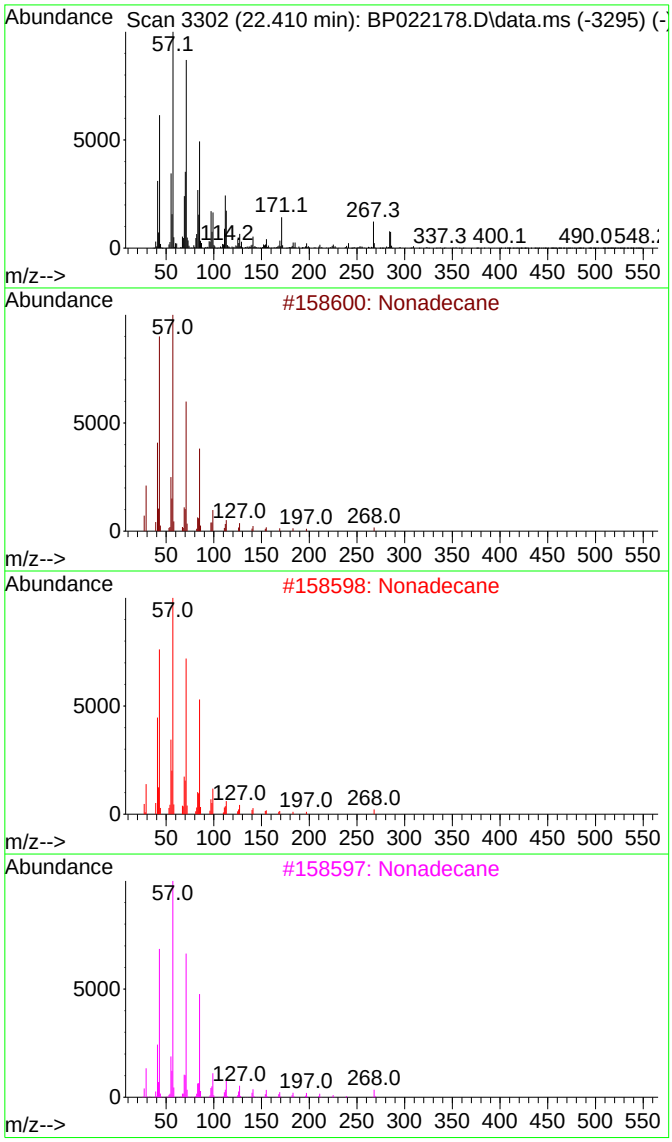
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

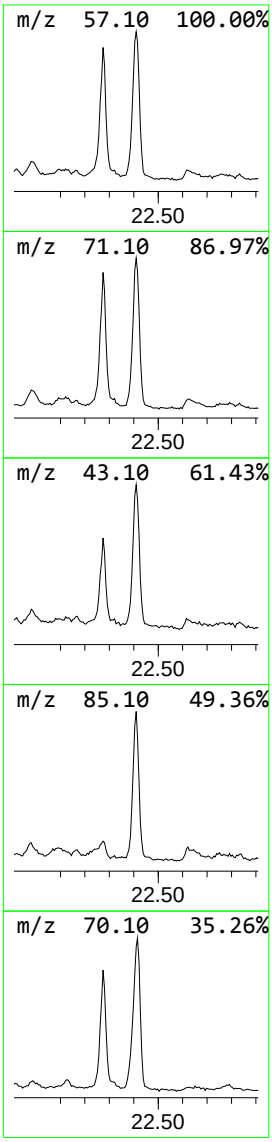
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.410	33.53 ng/ul	4634590	Chrysene-d12	21.533

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	95
2	Nonadecane	268	C19H40	000629-92-5	89
3	Nonadecane	268	C19H40	000629-92-5	76
4	Tetratetracontane	619	C44H90	007098-22-8	70
5	Octacosane	394	C28H58	000630-02-4	68





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

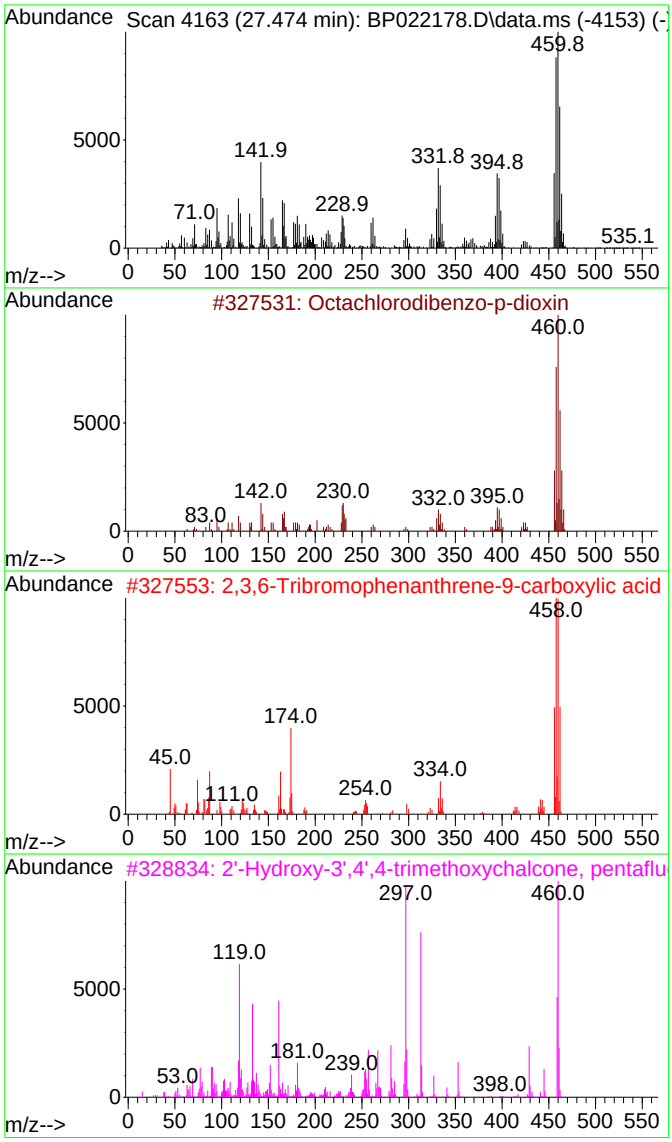
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

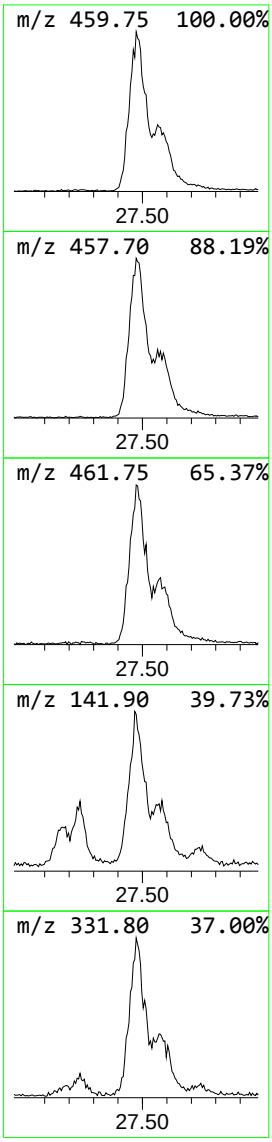
****

Peak Number 69 Octachlorodibenzo-p-dioxin Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.474	8.82 ng/ul	2330460	Perylene-d12	24.810

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	95
2	2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	42
3	2'-Hydroxy-3',4',4-trimethoxycha...	460	C21H17F5O6	1000453-83-5	9
4	Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	8
5	Moracin M, 3TMS	458	C23H34O4Si3	1000501-75-2	8





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022178.D
Acq On : 03 Oct 2024 02:18
Operator : RC/JU
Sample : P4017-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DDC70

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.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
(DEL) Alkane: S...	5.758	25.1	ng/ul	8399460	1	7.705	6692750	20.0
(DEL) Alkane: S...	6.222	8.0	ng/ul	2689750	1	7.705	6692750	20.0
(DEL) Alkane: S...	6.287	8.8	ng/ul	2934300	1	7.705	6692750	20.0
(DEL) Alkane: C...	6.369	11.5	ng/ul	3858630	1	7.705	6692750	20.0
(DEL) Alkane: S...	6.716	12.6	ng/ul	4206700	1	7.705	6692750	20.0
(DEL) Alkane: S...	6.769	14.8	ng/ul	4945360	1	7.705	6692750	20.0
Benzene, 1-ethy...	6.810	11.7	ng/ul	3919930	1	7.705	6692750	20.0
Benzene, 1-ethy...	7.099	11.8	ng/ul	3940130	1	7.705	6692750	20.0
2-Heptanone, 4,...	7.146	7.8	ng/ul	2595520	1	7.705	6692750	20.0
Carbonic acid, ...	7.352	84.5	ng/ul	28272900	1	7.705	6692750	20.0
(DEL) Alkane: S...	7.622	10.9	ng/ul	3663830	1	7.705	6692750	20.0
1-Hexanol, 2-et...	7.787	42.2	ng/ul	14121600	1	7.705	6692750	20.0
Mesitylene	7.822	10.8	ng/ul	3625450	1	7.705	6692750	20.0
(DEL) Alkane: C...	7.993	8.8	ng/ul	2930650	1	7.705	6692750	20.0
Cyclohexanone, ...	8.134	10.0	ng/ul	3335080	1	7.705	6692750	20.0
Benzene, 1-meth...	8.234	17.2	ng/ul	5768090	1	7.705	6692750	20.0
(DEL) Alkane: S...	8.293	7.7	ng/ul	2574630	1	7.705	6692750	20.0
Benzene, 2-ethy...	8.340	15.8	ng/ul	5273580	1	7.705	6692750	20.0
(DEL) Alkane: S...	8.463	15.1	ng/ul	5035070	1	7.705	6692750	20.0
Benzene, 1-meth...	8.499	11.7	ng/ul	3924670	1	7.705	6692750	20.0
Benzene, 4-ethy...	8.646	9.8	ng/ul	3280260	1	7.705	6692750	20.0
o-Cymene	8.699	12.0	ng/ul	4019810	1	7.705	6692750	20.0
(DEL) Alkane: S...	8.934	73.0	ng/ul	24411600	1	7.705	6692750	20.0
unknown-01	9.040	20.8	ng/ul	6944180	1	7.705	6692750	20.0
(DEL) Alkane: S...	9.616	2.7	ng/ul	2390280	2	10.463	17847200	20.0
unknown-02	9.998	4.8	ng/ul	4276640	2	10.463	17847200	20.0
unknown-03	10.840	10.4	ng/ul	9324640	2	10.463	17847200	20.0
(DEL) Alkane: S...	11.493	5.7	ng/ul	5113300	2	10.463	17847200	20.0
Benzene, 1,3,5-...	11.563	13.0	ng/ul	11561300	2	10.463	17847200	20.0
(DEL) Alkane: S...	11.722	3.4	ng/ul	3000790	2	10.463	17847200	20.0
Carbonic acid, ...	11.881	9.6	ng/ul	8582280	2	10.463	17847200	20.0
(DEL) Alkane: S...	11.922	2.9	ng/ul	2583910	2	10.463	17847200	20.0
Naphthalene, 1,...	13.481	4.0	ng/ul	6540660	3	14.310	32862600	20.0
Naphthalene, 2,...	13.628	3.9	ng/ul	6454700	3	14.310	32862600	20.0
(DEL) Alkane: S...	13.792	2.9	ng/ul	4684170	3	14.310	32862600	20.0
Carbonic acid, ...	14.234	20.0	ng/ul	32862600	3	14.310	32862600	20.0
1H-Indole, 4-(3...	14.463	8.7	ng/ul	14291700	3	14.310	32862600	20.0
Naphthalene, 1,...	15.098	10.0	ng/ul	16427600	3	14.310	32862600	20.0
(DEL) Alkane: S...	15.187	7.8	ng/ul	12772200	3	14.310	32862600	20.0
(DEL) Alkane: S...	15.598	2.7	ng/ul	4454330	3	14.310	32862600	20.0
unknown-04	15.975	14.0	ng/ul	4058370	4	17.145	5801300	20.0
(DEL) Alkane: S...	16.075	46.8	ng/ul	13561100	4	17.145	5801300	20.0
(DEL) Alkane: S...	16.933	28.7	ng/ul	8330520	4	17.145	5801300	20.0
2-Amino-5-isopr...	17.433	19.1	ng/ul	5526090	4	17.145	5801300	20.0
(DEL) Alkane: S...	17.639	30.1	ng/ul	8716900	4	17.145	5801300	20.0
2-(1,6,7-Ttrime...	17.722	9.6	ng/ul	2790250	4	17.145	5801300	20.0
Hexadecanoic ac...	17.822	8.8	ng/ul	2539940	4	17.145	5801300	20.0
n-Hexadecanoic ...	18.069	14.9	ng/ul	4316560	4	17.145	5801300	20.0
(DEL) Alkane: S...	18.357	23.7	ng/ul	6882260	4	17.145	5801300	20.0
(DEL) Alkane: S...	19.016	27.2	ng/ul	7882560	4	17.145	5801300	20.0
(DEL) Alkane: S...	20.174	34.0	ng/ul	4696110	5	21.533	2764580	20.0
(DEL) Alkane: S...	20.698	19.7	ng/ul	2727850	5	21.533	2764580	20.0

Carbonic acid, ...	21.216	64.9	ng/ul	8971370	5	21.533	2764580	20.0
(DEL) Alkane: S...	21.780	25.4	ng/ul	3512360	5	21.533	2764580	20.0
9-Octadecenoic ...	22.274	29.2	ng/ul	4034210	5	21.533	2764580	20.0
(DEL) Alkane: S...	22.410	33.5	ng/ul	4634590	5	21.533	2764580	20.0
Octachlorodiben...	27.474	8.8	ng/ul	2330460	6	24.810	5282750	20.0

Instrument :
BNA_P
ClientSampleId :
DDC70