LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110123\
Data File : BG059575.D
Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

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_G\Methods\SFAM-EPA-BG102723.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG059575.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.664	43	47	51	rVV6	10934	17277	1.46%	0.100%
2	3.728	57	58	63	rBV3	6021	6508	0.55%	0.037%
3	4.098	116	121	126	rBV3	25187	43122	3.65%	0.248%
4	4.198	135	138	139	rVV2	5034	6632	0.56%	0.038%
5	4.216	139	141	145	rVV3	9469	12128	1.03%	0.070%
6	4.245	145	146	150	rVV3	6097	6996	0.59%	0.040%
7	4.304	153	156	164	rVV3	18514	34053	2.88%	0.196%
8	4.439	177	179	184	rVB5	12981	17713	1.50%	0.102%
9	4.480	184	186	189	rBV2	6915	6782	0.57%	0.039%
10	5.368	333	337	340	rVV2	25714	42937	3.63%	0.247%
11	5.397	340	342	343	rVV2	8885	7388	0.62%	0.043%
12	5.456	348	352	354	rVV3	7826	10363	0.88%	0.060%
13	5.479	354	356	358	rVB3	8076	6624	0.56%	0.038%
14	5.614	373	379	385	rVB3	35449	69608	5.88%	0.401%
15	5.685	389	391	392	rBV2	9589	6744	0.57%	0.039%
16	5.908	426	429	430	rBV3	8654	9332	0.79%	0.054%
17	5.926	430	432	433	rVV2	11012	9706	0.82%	0.056%
18	5.955	436	437	442	rVB5	11089	8903	0.75%	0.051%
19	6.137	464	468	469	rBV	13069	12503	1.06%	0.072%
20	6.220	477	482	483	rBV4	6824	10994	0.93%	0.063%
21	6.243	483	486	488	rVV4	7918	9229	0.78%	0.053%
22	6.308	493	497	502	rVV5	15378	30146	2.55%	0.174%
23	6.378	504	509	511	rBV3	5950	7567	0.64%	0.044%
24	6.531	531	535	536	rBV3	9834	11449	0.97%	0.066%
25	6.731	565	569	570	rBV2	6305	6618	0.56%	0.038%
26	6.807	575	582	589	rBV7	31834	74168	6.27%	0.427%
27	6.925	596	602	610	rVB4	33618	70252	5.94%	0.405%
28	6.995	610	614	617	rBV5	5317	8295	0.70%	0.048%
29	7.107	629	633	634	rBV4	5683	6690	0.57%	0.039%
30	7.136	637	638	642	rVV5	9256	10204	0.86%	0.059%
31	7.165	642	643	647	rVV4	10201	7520	0.64%	0.043%
32	7.254	654	658	662	rVV4	25391	40943	3.46%	0.236%
33	7.295	662	665	672	rVB2	43583	80173	6.78%	0.462%
34	7.418	683	686	687	rBV3	17028	15474	1.31%	0.089%
35	7.465	687	694	703	rVV	210349	398345	33.67%	2.294%
36	7.577	710	713	716	rBV5	5539	6549	0.55%	0.038%

LSC Area Percent Report

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Misc :
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Instrument :
BNA_G
ClientSampleId :
EOAQ3

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_G\Methods\SFAM-EPA-BG102723.MA.M
Title : SVOA CALIBRATION

37	7.677	724	730	739	rBV2	140186	261865	22.14%	1.508%
38	7.782	745	748	753	rBV5	16568	22224	1.88%	0.128%
39	7.871	757	763	777	rVB2	200255	392494	33.18%	2.261%
40	7.976	777	781	785	rVB4	27226	37416	3.16%	0.216%
41	8.017	785	788	790	rBV3	9764	11613	0.98%	0.067%
42	8.247	820	827	836	rVB5	44409	108819	9.20%	0.627%
43	8.358	839	846	857	rBV	315024	544189	46.00%	3.134%
44	8.464	857	864	868	rBV9	22024	46241	3.91%	0.266%
45	8.511	868	872	873	rBV4	6658	8979	0.76%	0.052%
46	8.593	882	886	891	rBV7	20373	40218	3.40%	0.232%
47	8.717	902	907	918	rBV2	48953	122122	10.32%	0.703%
48	8.816	918	924	930	rVV7	35299	81131	6.86%	0.467%
49	8.875	930	934	939	rVB6	38856	73088	6.18%	0.421%
50	8.981	942	952	956	rBV9	45174	118061	9.98%	0.680%
51	9.040	956	962	973	rVB2	225061	426827	36.08%	2.458%
52	9.198	982	989	996	rVB3	70477	163617	13.83%	0.942%
53	9.292	1000	1005	1006	rBV3	13346	22075	1.87%	0.127%
54	9.339	1011	1013	1016	rVB4	10844	10316	0.87%	0.059%
55	9.386	1016	1021	1024	rBV5	13144	18678	1.58%	0.108%
56	9.439	1024	1030	1039	rVV4	86141	181227	15.32%	1.044%
57	9.551	1040	1049	1055	rVB2	184363	394233	33.33%	2.271%
58	9.686	1067	1072	1079	rVB9	32579	76197	6.44%	0.439%
59	9.774	1079	1087	1091	rBV8	21837	62626	5.29%	0.361%
60	9.892	1103	1107	1111	rVB6	21896	30192	2.55%	0.174%
61	9.933	1112	1114	1117	rVV4	23169	32186	2.72%	0.185%
62	9.974	1117	1121	1126	rVB3	54299	97470	8.24%	0.561%
63	10.033	1129	1131	1132	rVV2	17912	14529	1.23%	0.084%
64	10.068	1135	1137	1144	rVB5	35355	55687	4.71%	0.321%
65	10.156	1150	1152	1154	rBV3	10932	12746	1.08%	0.073%
66	10.268	1161	1171	1178	rVB5	170491	405128	34.25%	2.333%
67	10.350	1178	1185	1193	rBV9	79362	206779	17.48%	1.191%
68	10.415	1193	1196	1197	rBV	33480	32602	2.76%	0.188%
69	10.562	1211	1221	1227	rBV3	110787	240019	20.29%	1.382%
70	10.738	1245	1251	1255	rVB8	27600	56400	4.77%	0.325%
71	10.797	1255	1261	1268	rBV2	252734	483979	40.91%	2.788%
72	10.849	1268	1270	1276	rVB3	41597	58881	4.98%	0.339%
73	10.908	1278	1280	1282	rBV3	10690	7466	0.63%	0.043%
74	11.084	1306	1310	1313	rVB6	25196	38818	3.28%	0.224%
75	11.208	1324	1331	1337	rBV	418401	767749	64.90%	4.422%
76	11.267	1337	1341	1349	rVB4	112162	231450	19.57%	1.333%

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Misc :
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Instrument :
BNA_G
ClientSampleId :
EOAQ3

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

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Title : SVOA CALIBRATION

77	11.349	1349	1355	1366	rVB3	142775	345893	29.24%	1.992%
78	11.760	1423	1425	1428	rBV4	20196	21109	1.78%	0.122%
79	11.836	1433	1438	1449	rVB8	94068	237900	20.11%	1.370%
80	11.936	1451	1455	1460	rBV8	45747	80593	6.81%	0.464%
81	12.119	1483	1486	1491	rVB	133295	186390	15.76%	1.074%
82	12.483	1545	1548	1552	rVB2	82880	114708	9.70%	0.661%
83	12.730	1584	1590	1594	rBV8	55975	117579	9.94%	0.677%
84	12.835	1603	1608	1613	rBV	142025	245927	20.79%	1.417%
85	13.047	1640	1644	1650	rBV3	164061	303968	25.70%	1.751%
86	13.135	1657	1659	1663	rBV5	43086	58546	4.95%	0.337%
87	13.399	1701	1704	1708	rVB5	54047	67275	5.69%	0.387%
88	13.452	1708	1713	1717	rBV3	151809	228032	19.28%	1.313%
89	14.140	1825	1830	1831	rBV3	96687	132957	11.24%	0.766%
90	14.304	1854	1858	1862	rVB3	144717	217894	18.42%	1.255%
91	14.381	1865	1871	1880	rVB2	668587	1182922	100.00%	6.814%
92	14.686	1918	1923	1929	rVB2	536085	866621	73.26%	4.992%
93	14.986	1969	1974	1981	rVB	571189	863547	73.00%	4.974%
94	15.150	1997	2002	2006	rBV3	212909	408606	34.54%	2.354%
95	15.973	2137	2142	2150	rVB	623938	994281	84.05%	5.727%
96	16.149	2168	2172	2178	rVB3	342665	546348	46.19%	3.147%
97	16.631	2250	2254	2258	rBV	454645	650332	54.98%	3.746%
98	17.453	2390	2394	2400	rBV	415812	600494	50.76%	3.459%
99	17.735	2438	2442	2447	rBV	617463	836113	70.68%	4.816%
100	17.835	2455	2459	2464	rVB	647717	927115	78.37%	5.340%

Sum of corrected areas: 17361422

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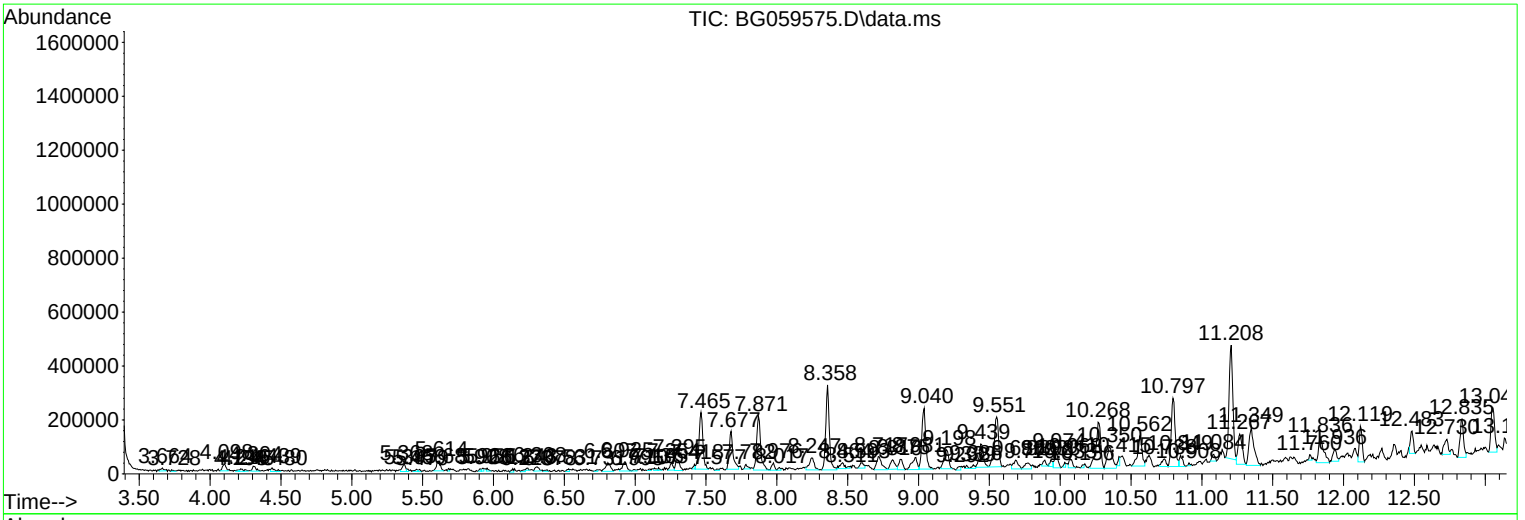
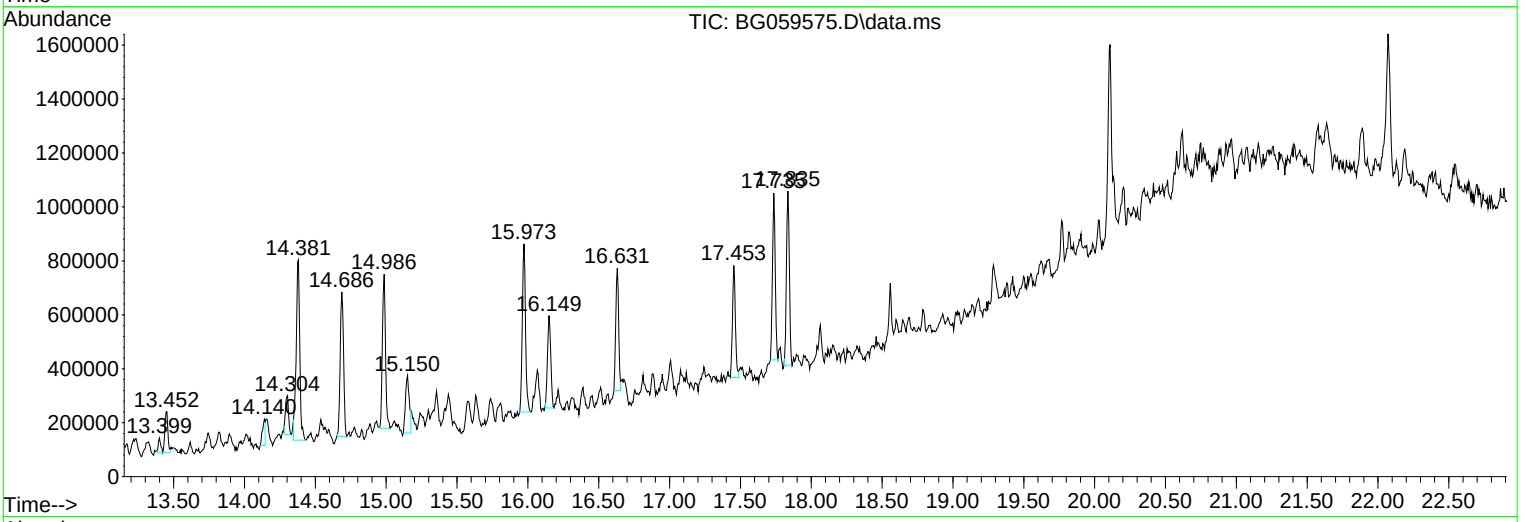
Misc :
ALS Vial : 4 Sample Multiplier: 1

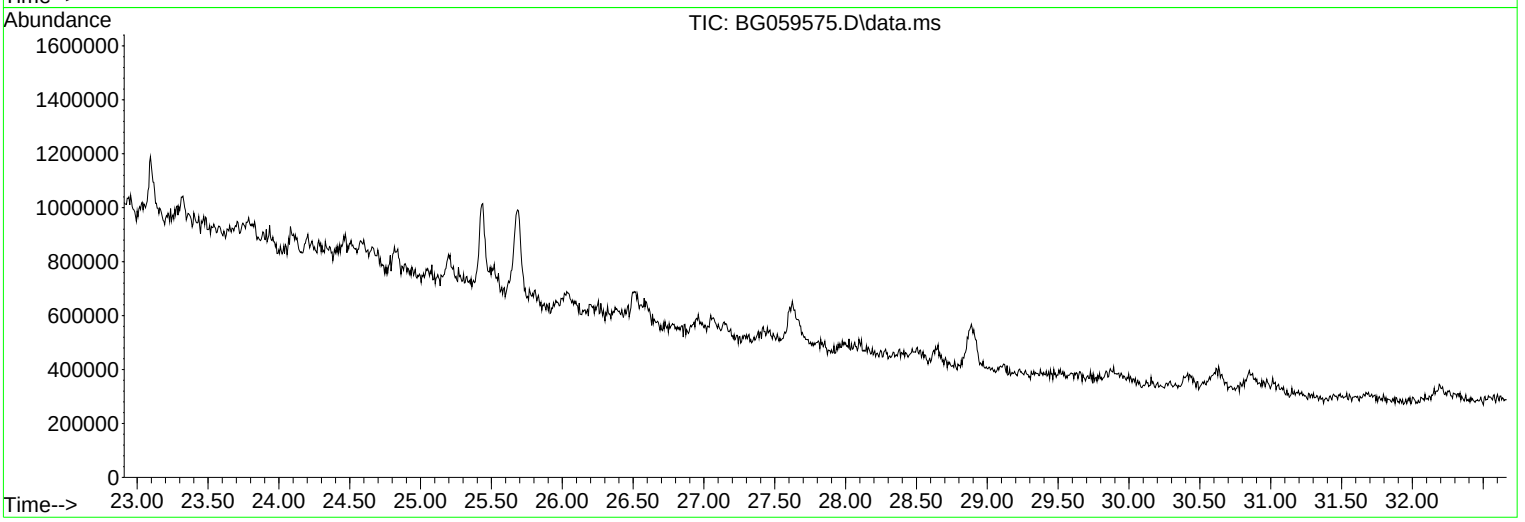
Instrument :
BNA_G
ClientSampleId :
EOAQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110123\
Data File : BG059575.D
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

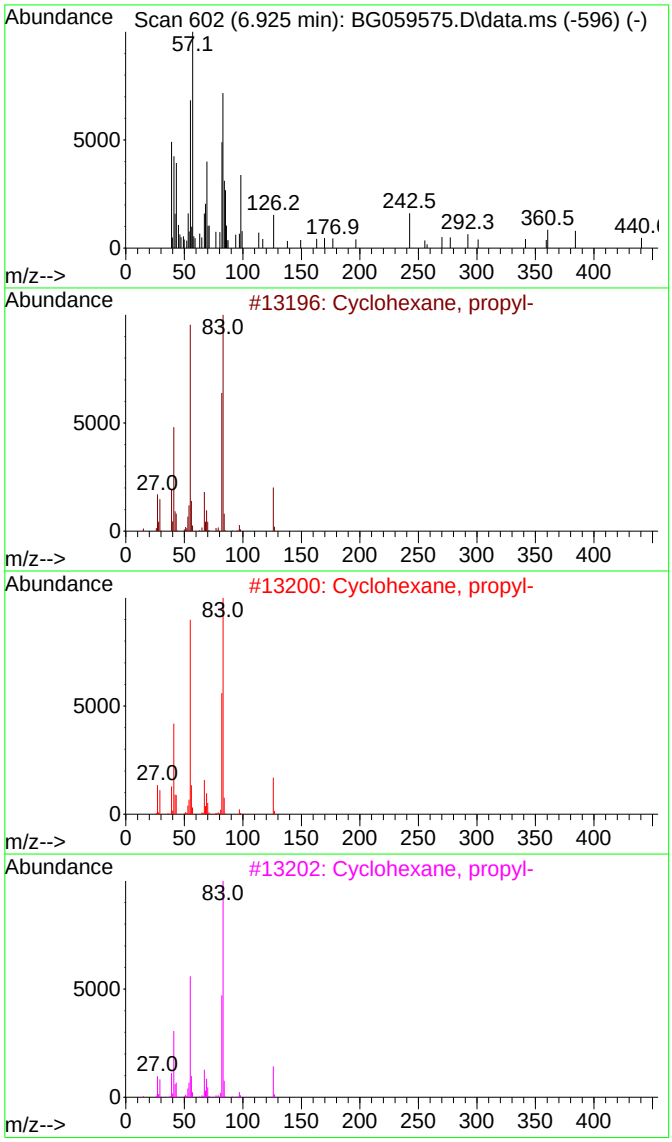
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

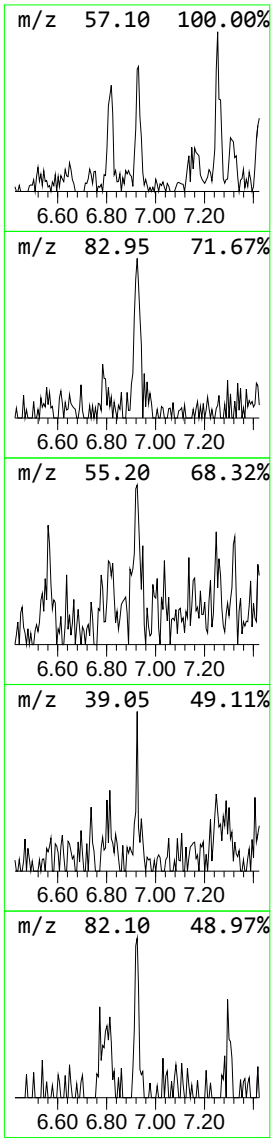
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Cyclic6.925 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.925	2.58 ng/ul	70252	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	43
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	43
3	Cyclohexane, propyl-	126	C9H18	001678-92-8	43
4	Cyclohexane, propyl-	126	C9H18	001678-92-8	43
5	Cyclohexane, propyl-	126	C9H18	001678-92-8	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110123\
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Sample : 04922-01
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

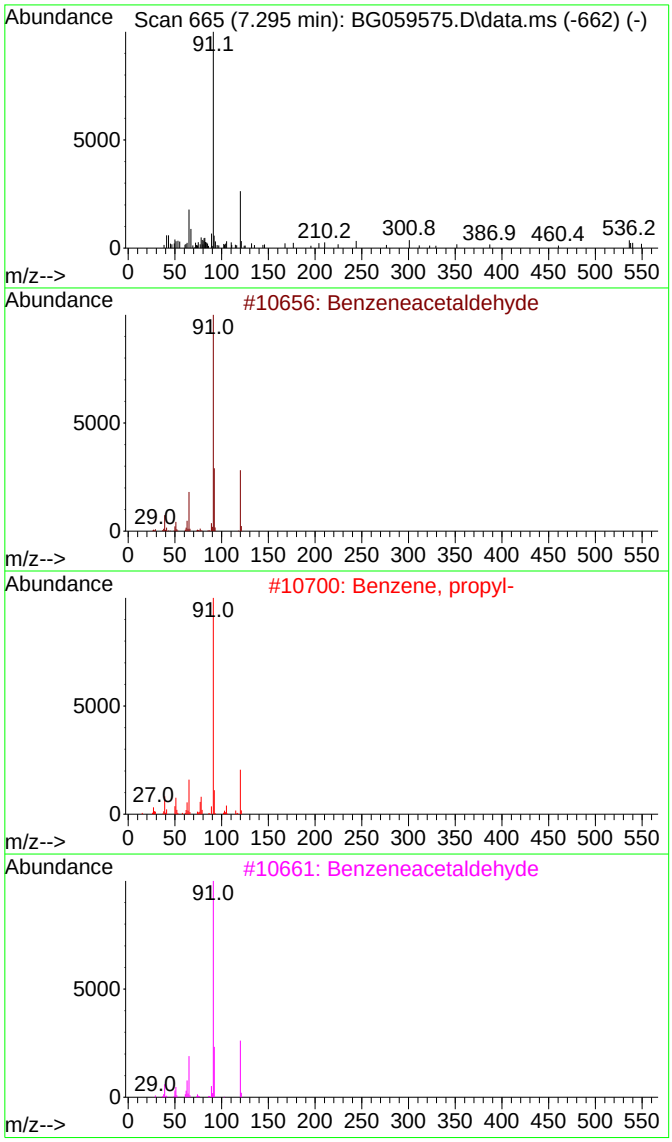
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

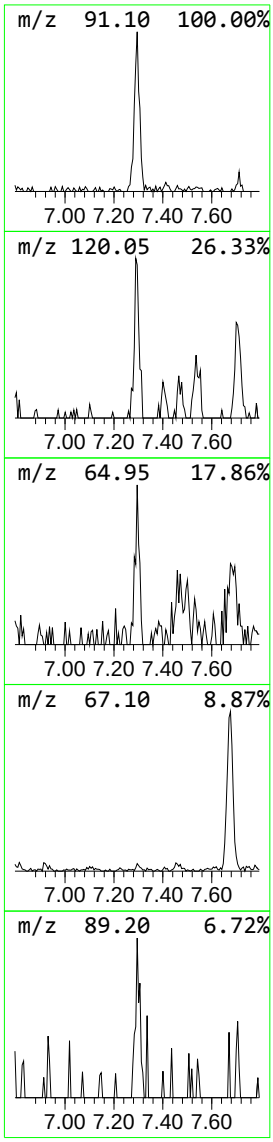
****

Peak Number 4 unknown-01 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.295	2.95 ng/ul	80173	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetaldehyde	120	C8H8O	000122-78-1	49
2	Benzene, propyl-	120	C9H12	000103-65-1	47
3	Benzeneacetaldehyde	120	C8H8O	000122-78-1	47
4	N-Isobutyl(phenyl)methanesulfona...	227	C11H17NO2S	144615-94-1	47
5	Benzeneacetaldehyde	120	C8H8O	000122-78-1	47





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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

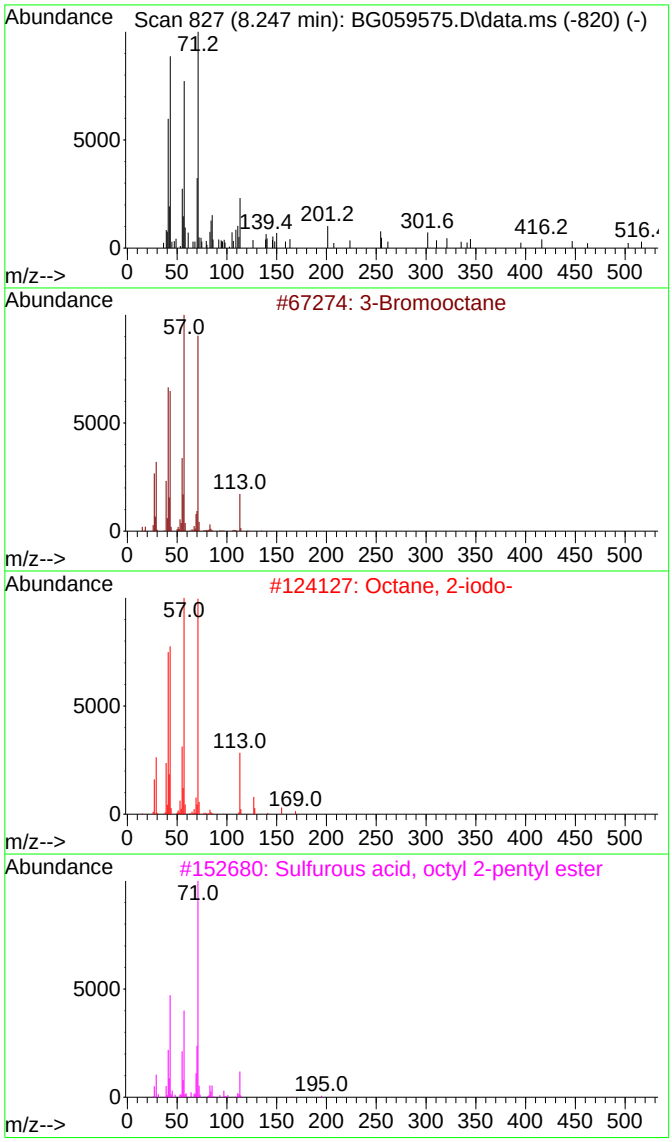
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

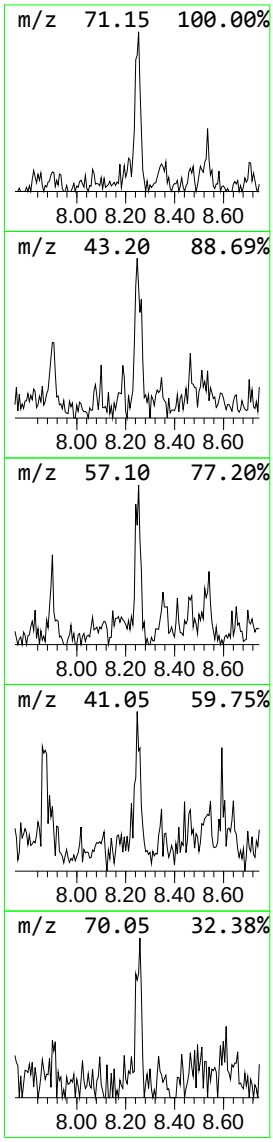
****

Peak Number 5 3-Bromooctane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.247	4.00 ng/ul	108819	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Bromooctane	192	C8H17Br	000999-64-4	59
2	Octane, 2-iodo-	240	C8H17I	000557-36-8	59
3	Sulfurous acid, octyl 2-pentyl e...	264	C13H28O3S	1000309-15-7	59
4	Decane, 4-methyl-	156	C11H24	002847-72-5	50
5	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	50





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Instrument :
BNA_G
ClientSampleId :
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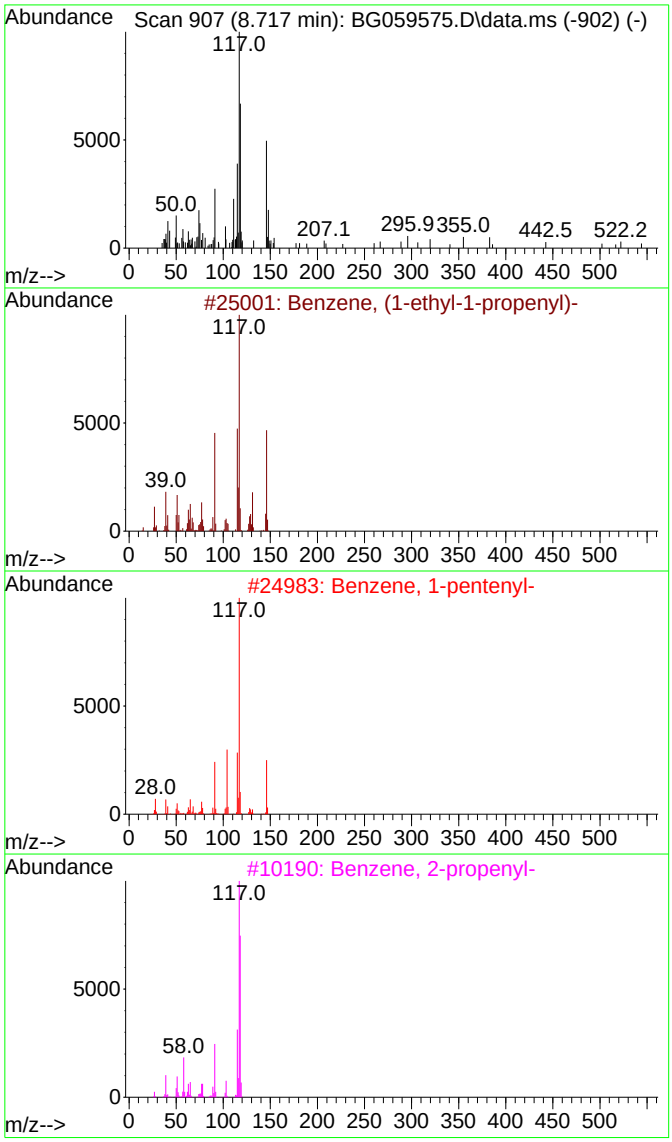
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

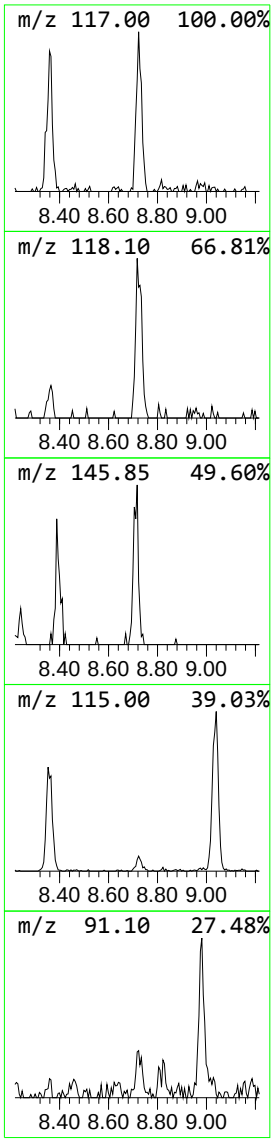
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, (1-ethyl-1-propenyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.717	4.49 ng/ul	122122	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	53
2	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	49
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	43
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	43
5	Indane	118	C9H10	000496-11-7	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110123\
Data File : BG059575.D
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Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

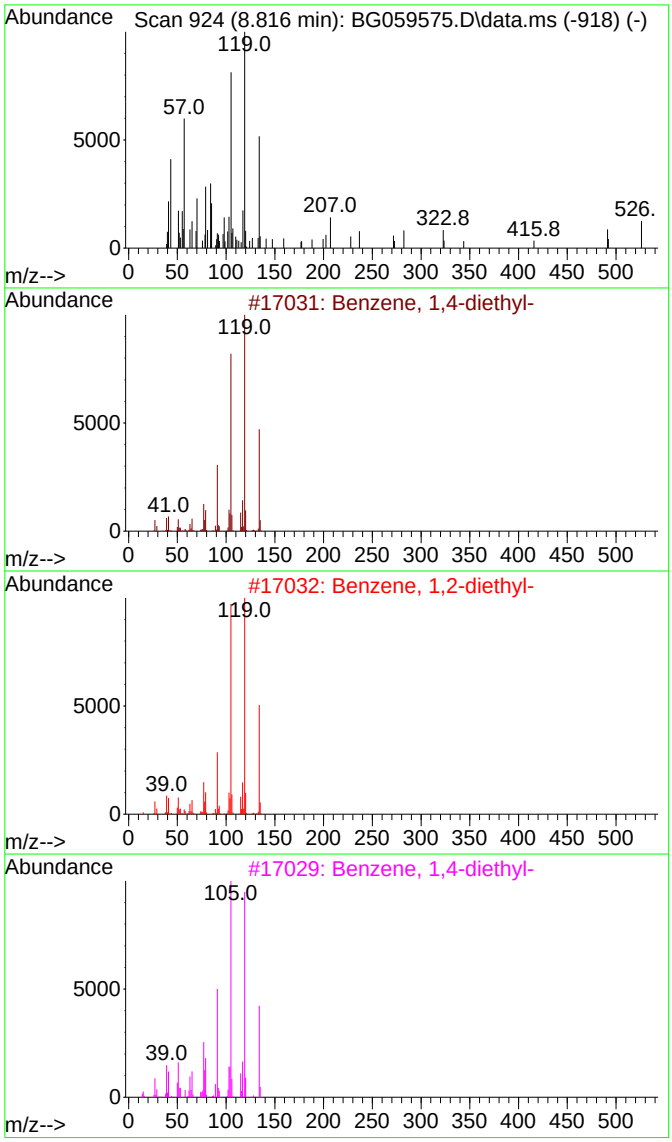
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

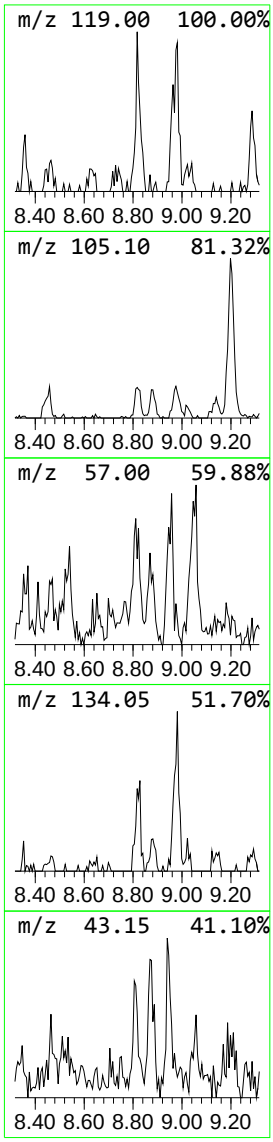
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.816	2.98 ng/ul	81131	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	53
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	49
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	49
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	49
5	Ethanone, 1-(3-methylphenyl)-	134	C9H10O	000585-74-0	49





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

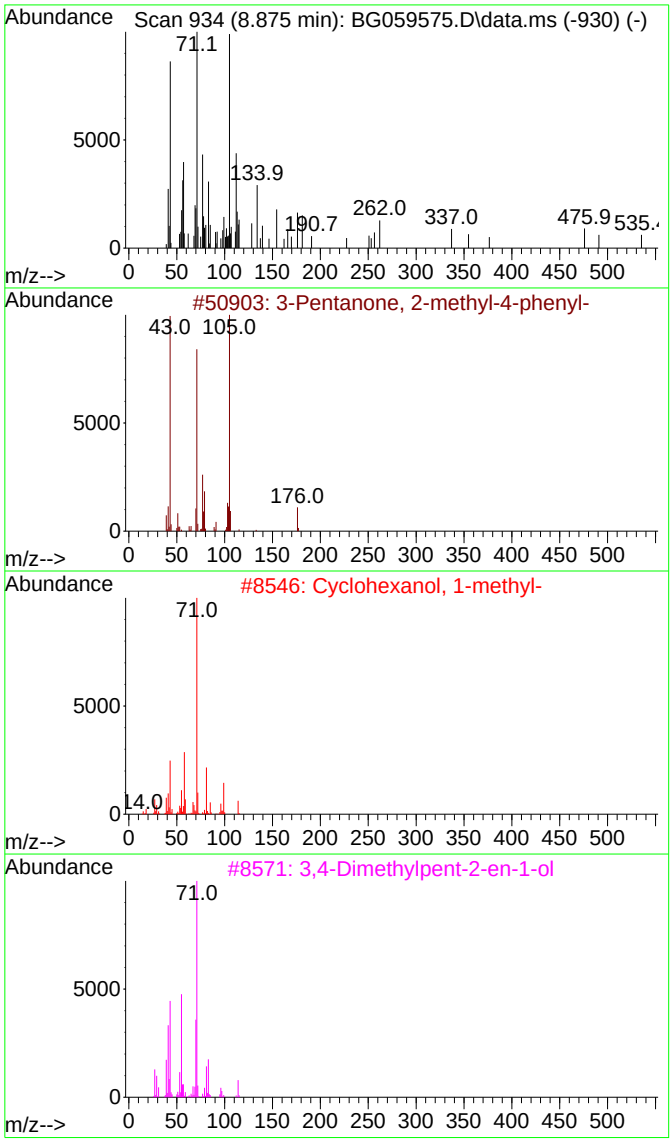
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

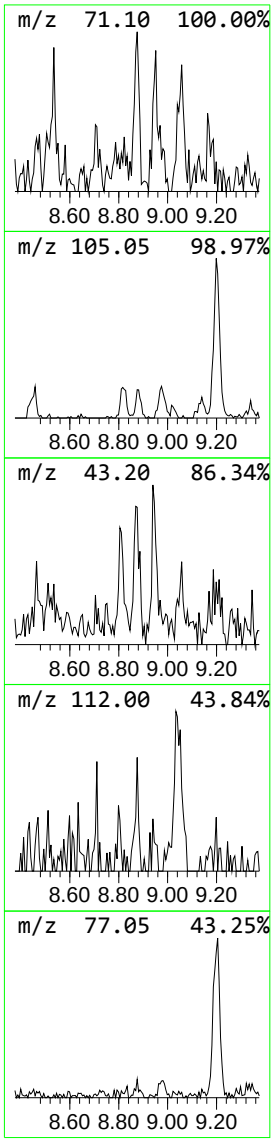
****

Peak Number 8 unknown-02 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.875	2.69 ng/ul	73088	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanone, 2-methyl-4-phenyl-	176	C12H16O	020474-49-1	38
2	Cyclohexanol, 1-methyl-	114	C7H14O	000590-67-0	30
3	3,4-Dimethylpent-2-en-1-ol	114	C7H14O	1000195-06-9	27
4	2-Hexen-1-ol, 3-methyl-, (E)-	114	C7H14O	030801-96-8	27
5	2,4-Pentanedione, 3-butyl-	156	C9H16O2	001540-36-9	27





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Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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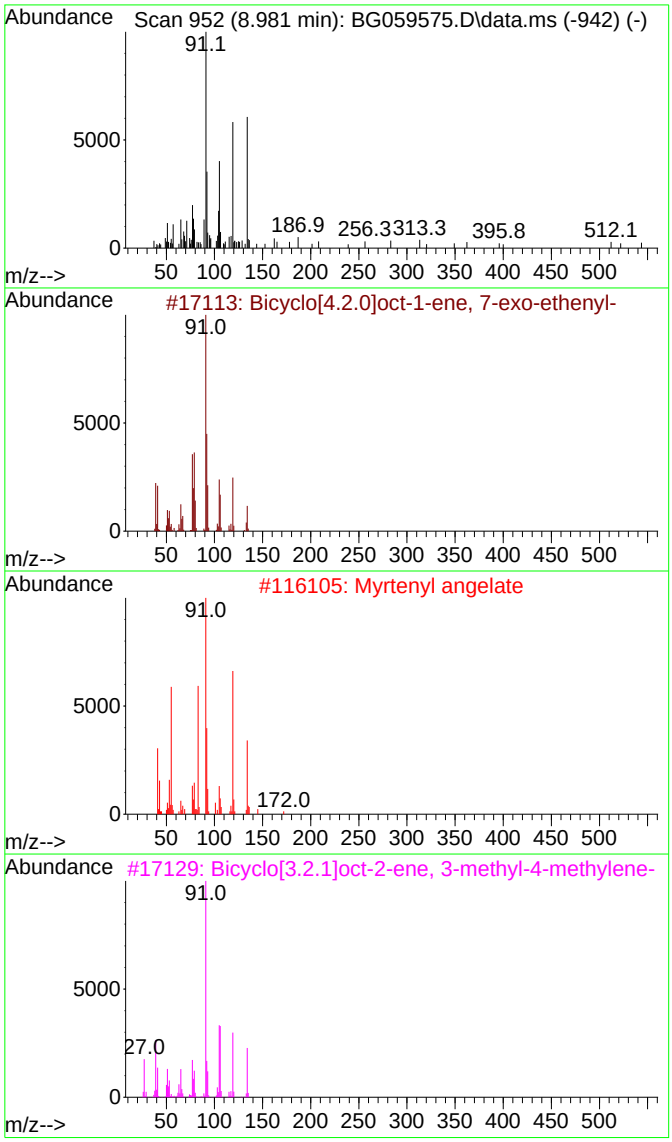
Quant Title : SVOA CALIBRATION

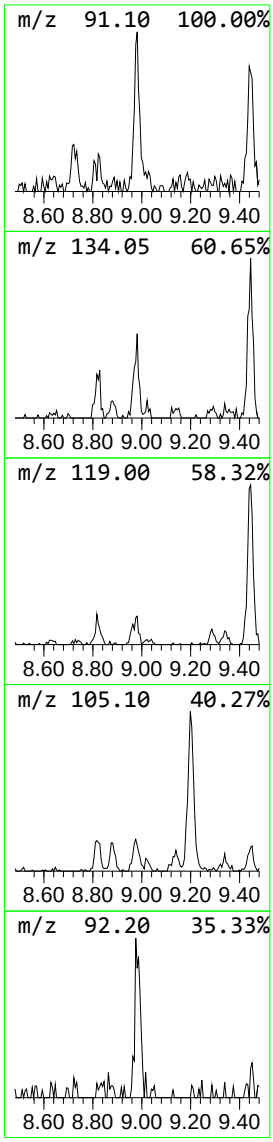
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Bicyclo[4.2.0]oct-1-ene, 7-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.981	4.34 ng/ul	118061	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]oct-1-ene, 7-exo-e...	134	C10H14	1000142-18-2	59
2	Myrtenyl angelate	234	C15H22O2	138530-45-7	50
3	Bicyclo[3.2.1]oct-2-ene, 3-methy...	134	C10H14	049826-53-1	50
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	50
5	1,2,3,4,5,8-Hexahydronaphthalene	134	C10H14	036231-13-7	50





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Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

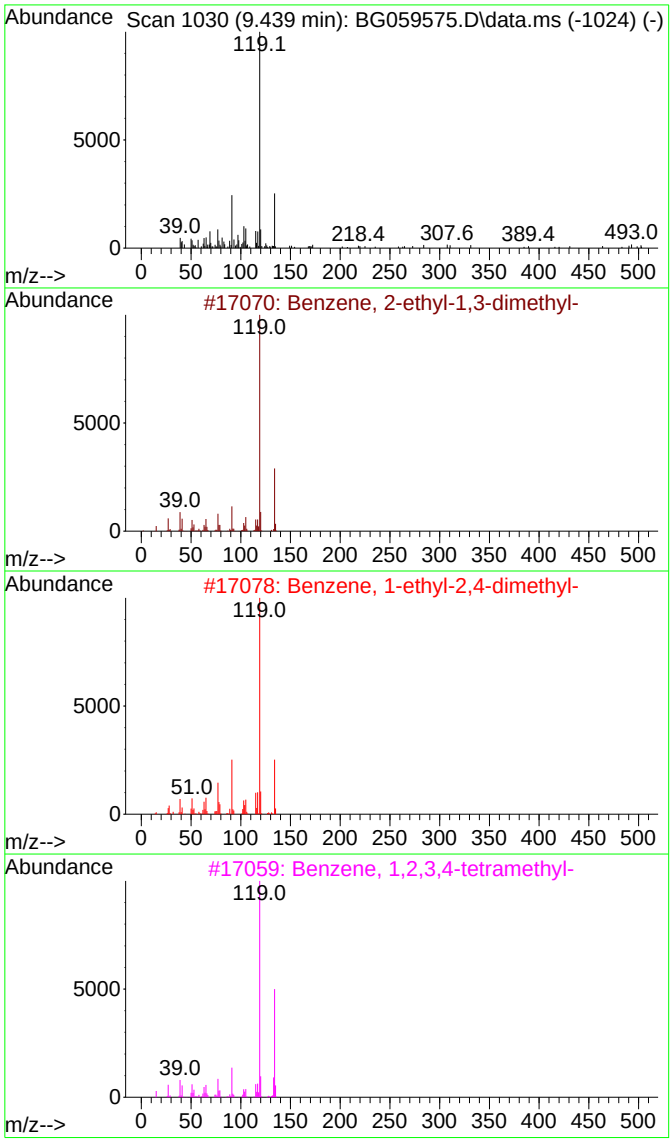
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

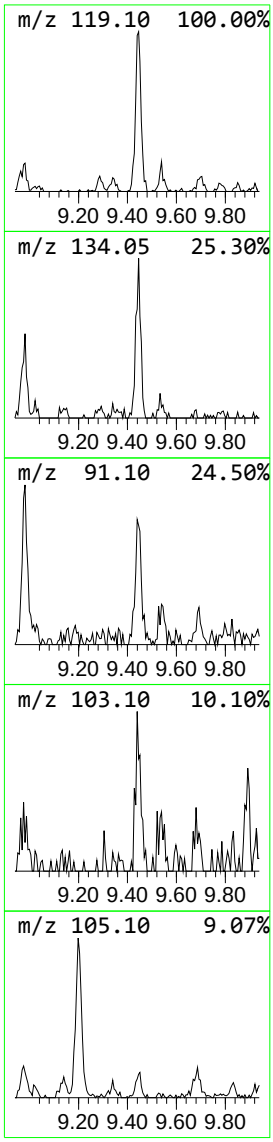
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.439	6.66 ng/ul	181227	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
2	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	87
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	87
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87





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Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

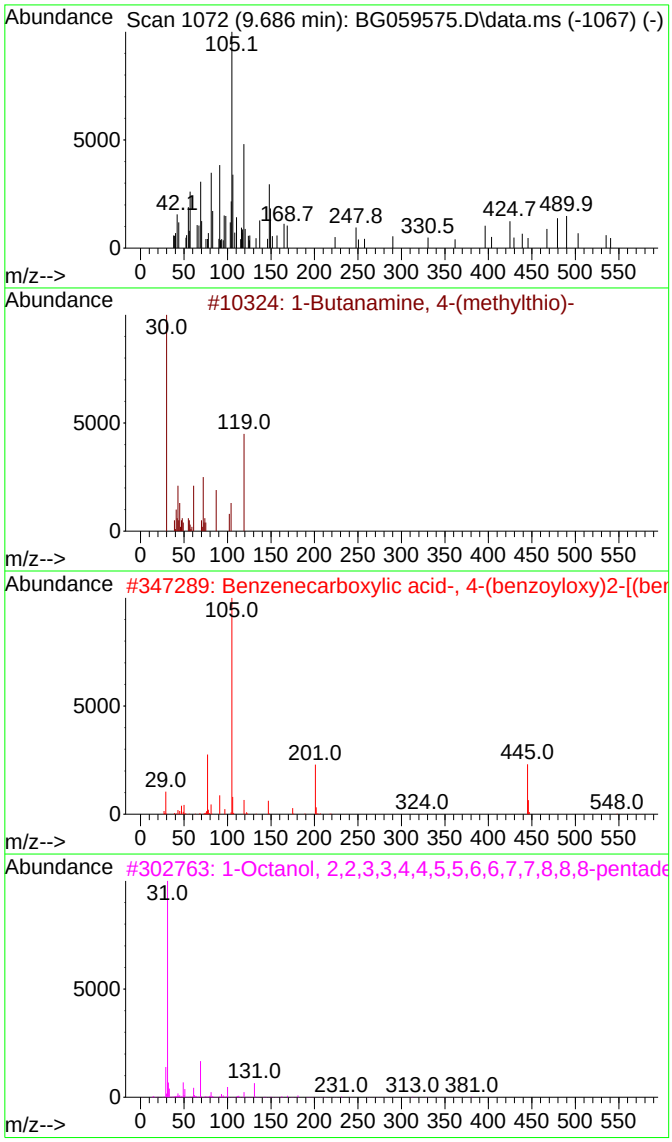
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

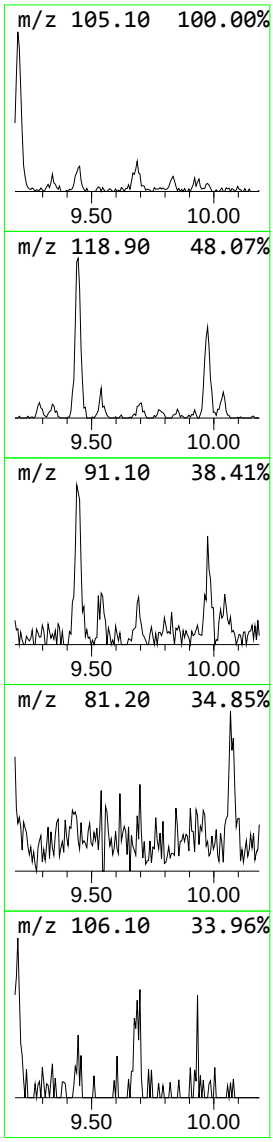
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 unknown-03 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.686	2.80 ng/ul	76197	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanamine, 4-(methylthio)-	119	C5H13NS	055021-77-7	3
2	Benzenecarboxylic acid-, 4-(benz...	620	C34H36O11	1000188-36-6	1
3	1-Octanol, 2,2,3,3,4,4,5,5,6,6,7...	400	C8H3F15O	000307-30-2	1
4	Propanoic acid, pentafluoro-, et...	192	C5H5F5O2	000426-65-3	1
5	3-(Perfluorohexyl)propanol-1	378	C9H7F13O	080806-68-4	1





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Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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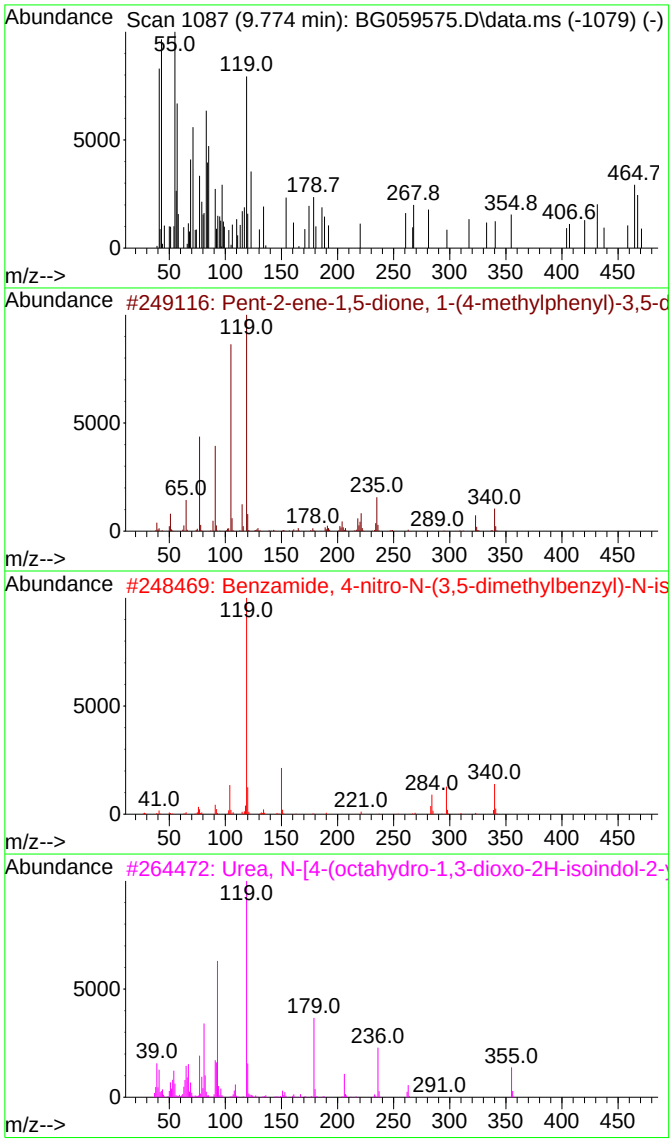
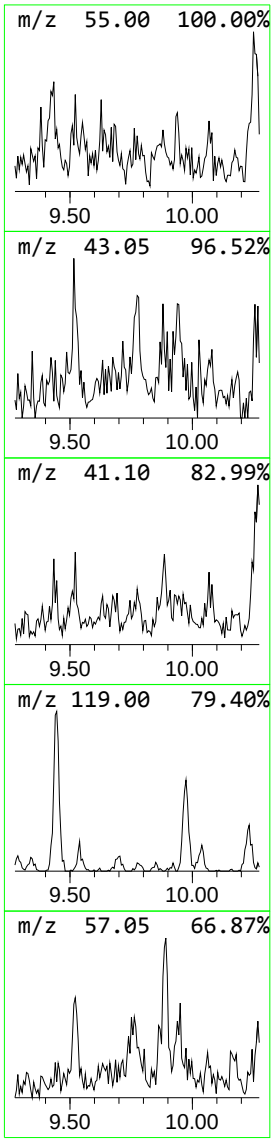
Peak Number 12 unknown-04 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.774	2.30 ng/ul	62626	1,4-Dichlorobenzene-d4	8.358

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pent-2-ene-1,5-dione, 1-(4-methy...	340	C24H20O2	063688-28-8	22
2	Benzamide, 4-nitro-N-(3,5-dimeth...	340	C20H24N2O3	1000491-64-2	12
3	Urea, N-[4-(octahydro-1,3-dioxo-...	355	C17H17N5O4	1000349-93-2	10
4	3-t-Butyl-1-p-tolyl-5-(4-trifluo...	390	C23H25F3O2	1000187-44-1	10
5	Sarcosine, N-(3-methylbenzoyl)-,...	333	C20H31NO3	1000321-15-9	9





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Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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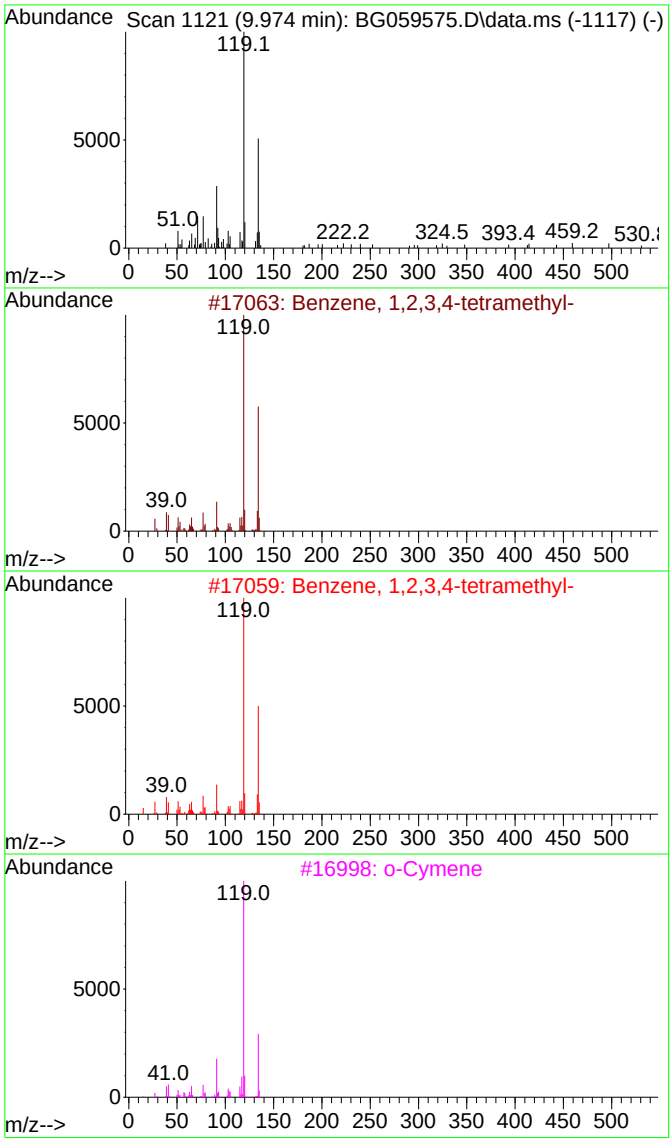
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

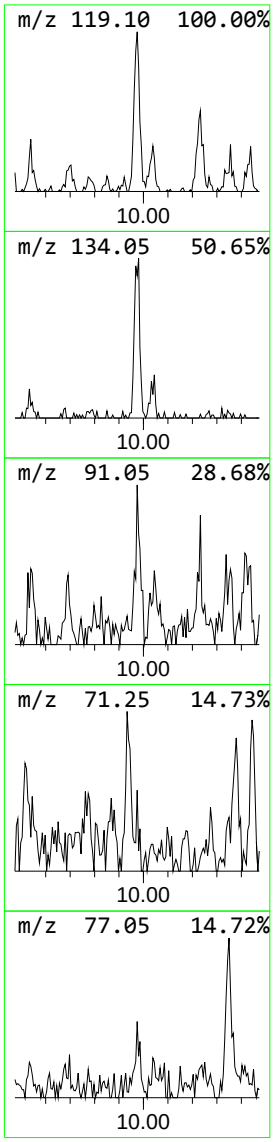
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.974	2.54 ng/ul	97470	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
3	o-Cymene	134	C10H14	000527-84-4	83
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	83
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	83





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Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

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

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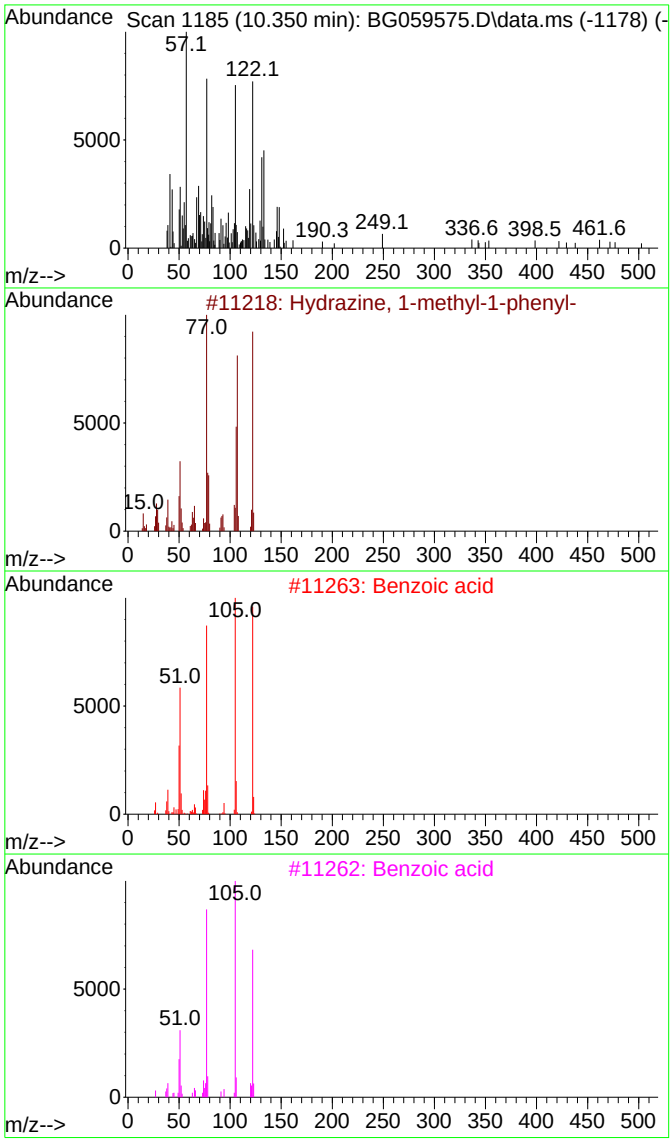
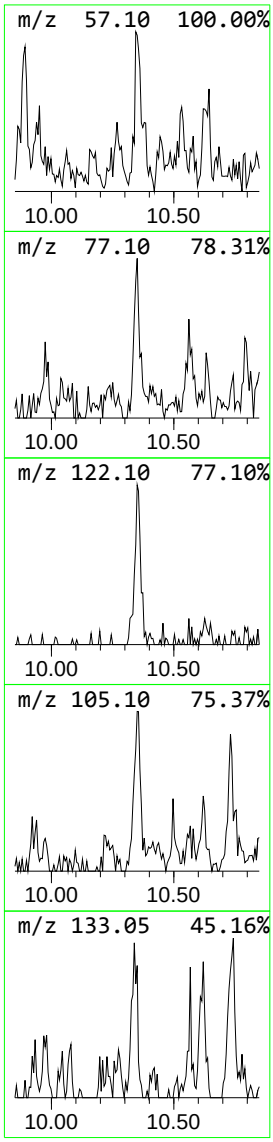
Peak Number 14 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.350	5.39 ng/ul	206779	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrazine, 1-methyl-1-phenyl-	122	C7H10N2	000618-40-6	49
2	Benzoic acid	122	C7H6O2	000065-85-0	49
3	Benzoic acid	122	C7H6O2	000065-85-0	43
4	(+)-Dibenzoyl-L-tartaric acid an...	340	C18H12O7	064339-95-3	43
5	Heptanediamide, N,N'-di-benzoyloxy-	398	C21H22N2O6	1000253-26-4	43





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Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

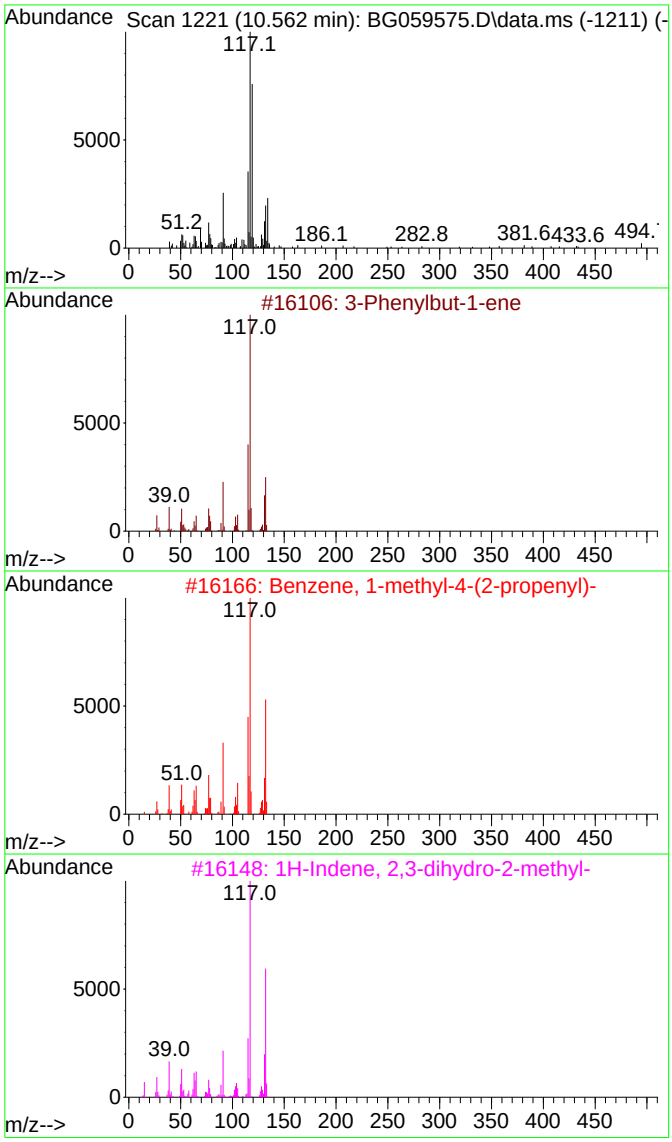
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

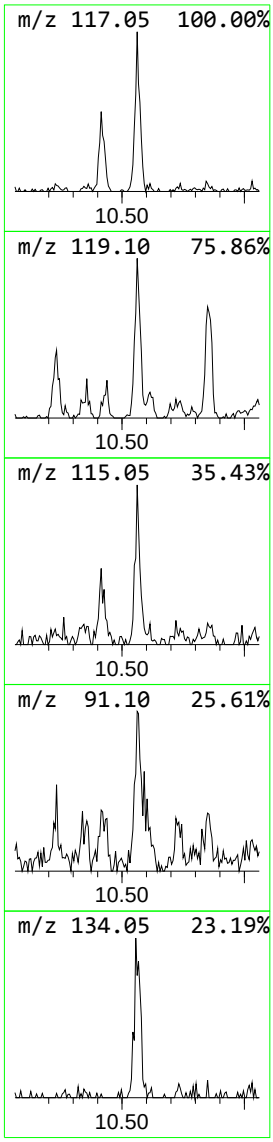
****

Peak Number 15 unknown-06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.562	6.25 ng/ul	240019	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	49
2	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	46
3	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	46
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	43
5	Indan, 1-methyl-	132	C10H12	000767-58-8	43





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Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

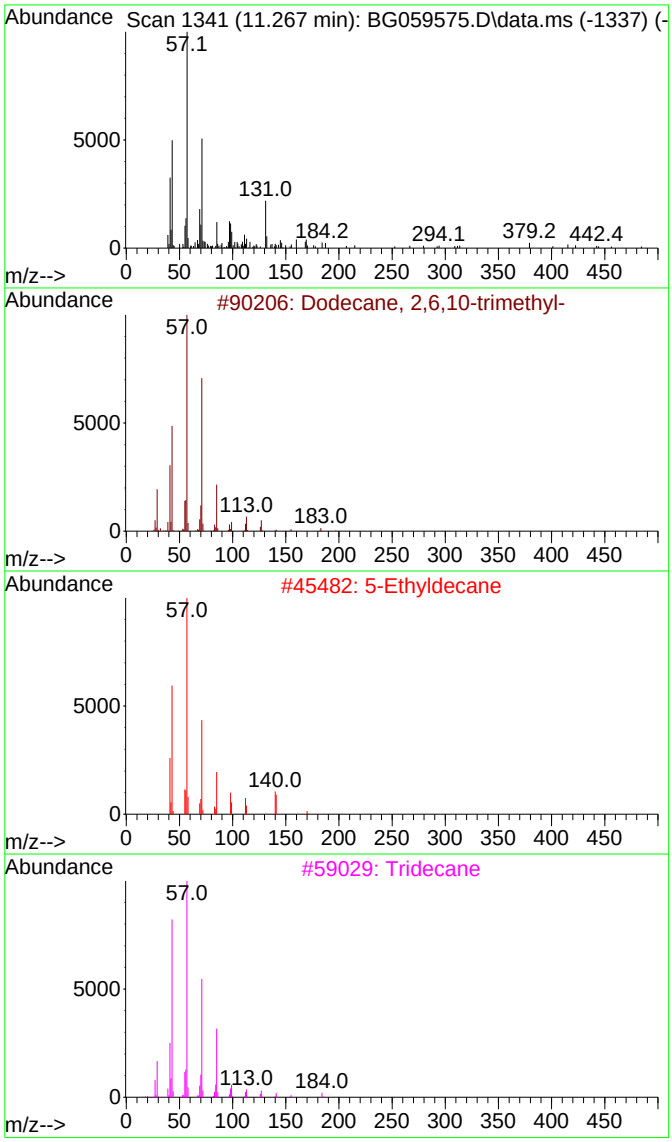
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

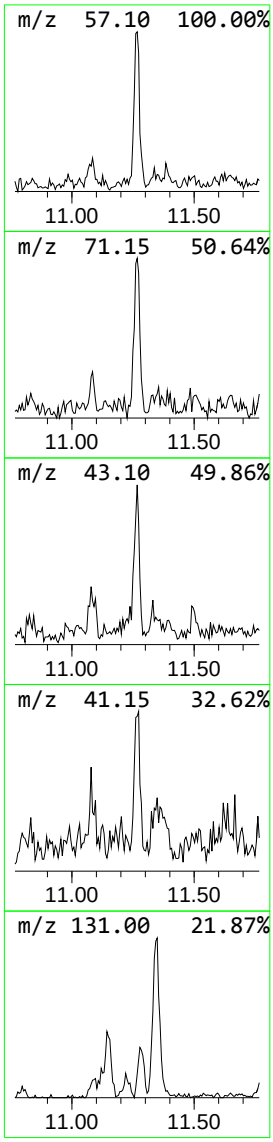
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.267	6.03 ng/ul	231450	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	47
2	5-Ethyldecane	170	C12H26	017302-36-2	46
3	Tridecane	184	C13H28	000629-50-5	46
4	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	46
5	Dodecane	170	C12H26	000112-40-3	43





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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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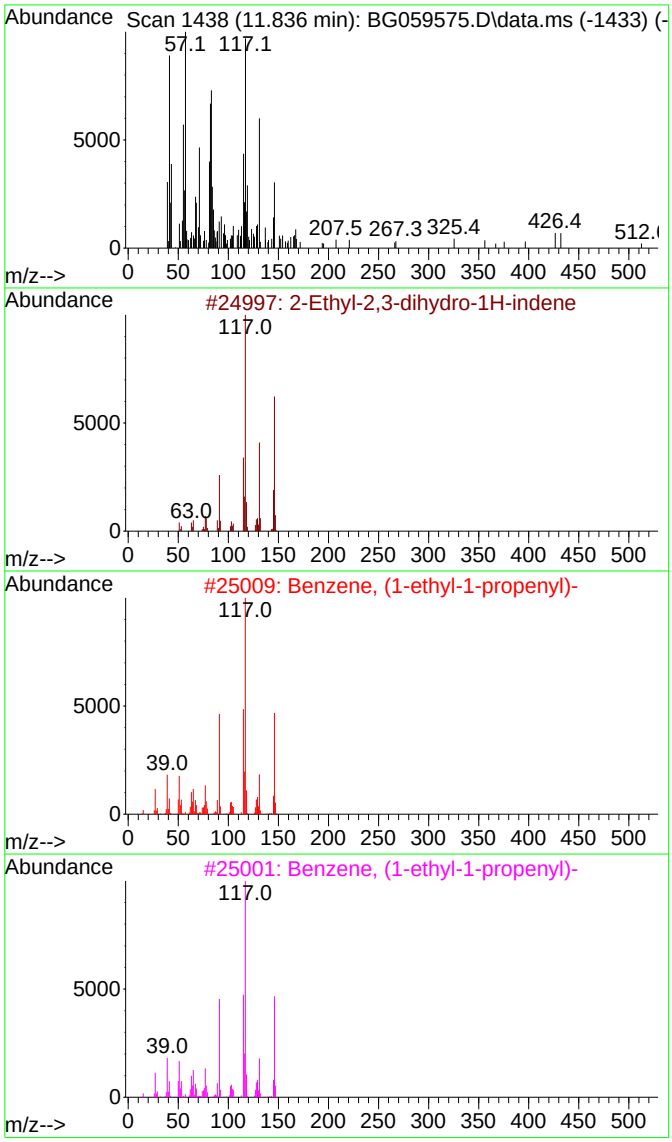
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

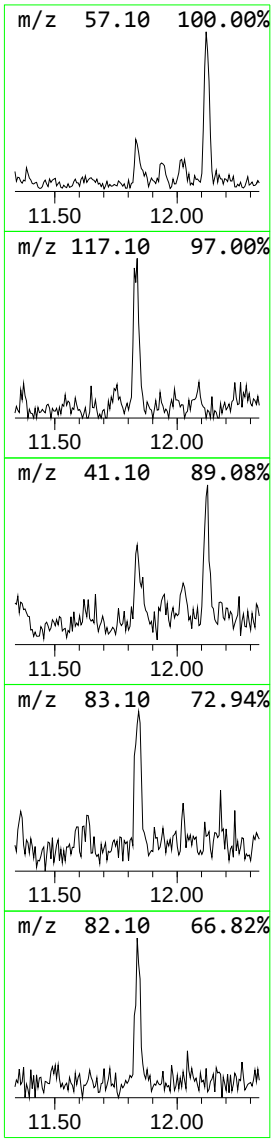
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 2-Ethyl-2,3-dihydro-1H-indene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.836	6.20 ng/ul	237900	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	81
2	Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	50
3	Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	50
4	trans-1-Phenyl-1-pentene	146	C11H14	016002-93-0	47
5	1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	45





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Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

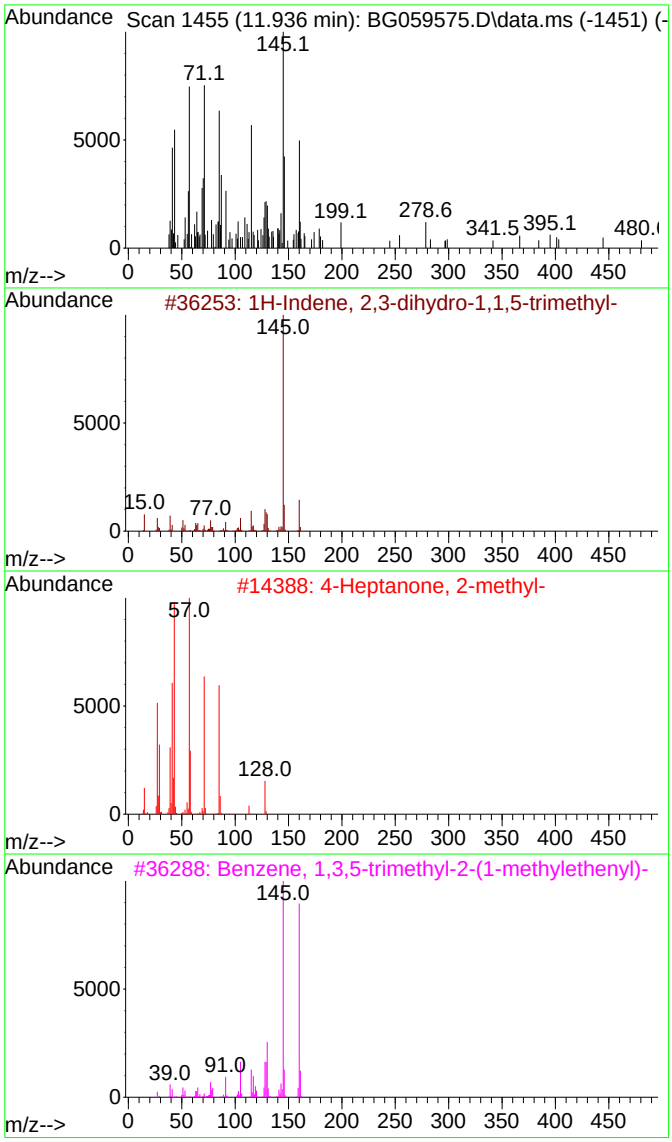
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

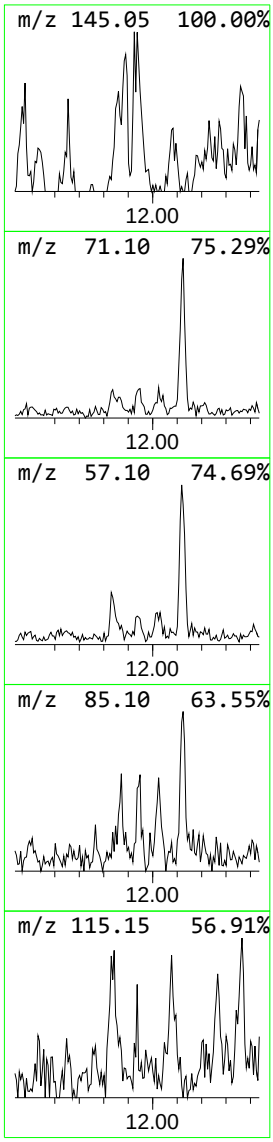
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 unknown-07 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.936	2.10 ng/ul	80593	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	38
2	4-Heptanone, 2-methyl-	128	C8H16O	000626-33-5	30
3	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	30
4	1H-Benzimidazole, 1-ethyl-2-methyl-	160	C10H12N2	005805-76-5	25
5	Heneicosane	296	C21H44	000629-94-7	25





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Data File : BG059575.D
Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

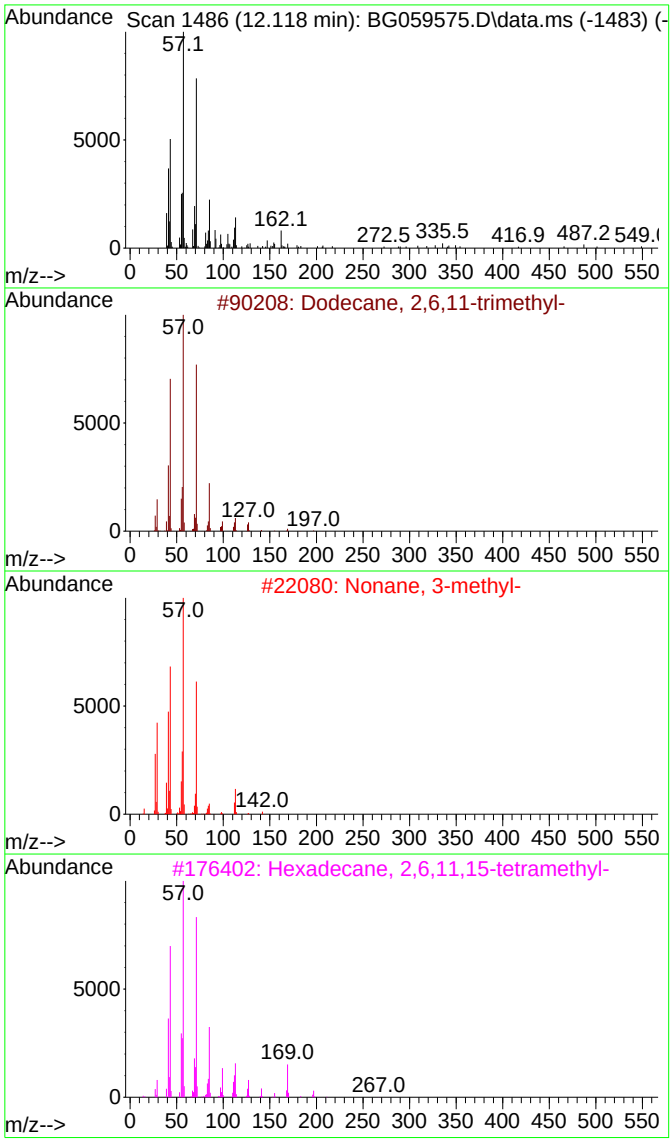
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

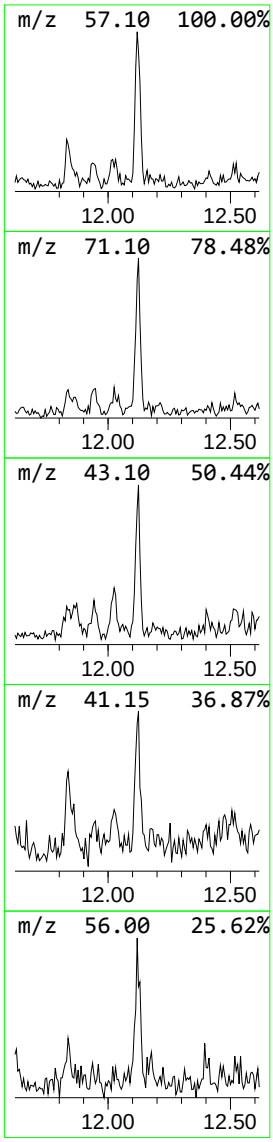
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.118	4.86 ng/ul	186390	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	90
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	83
3	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	83
4	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	80
5	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	80





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

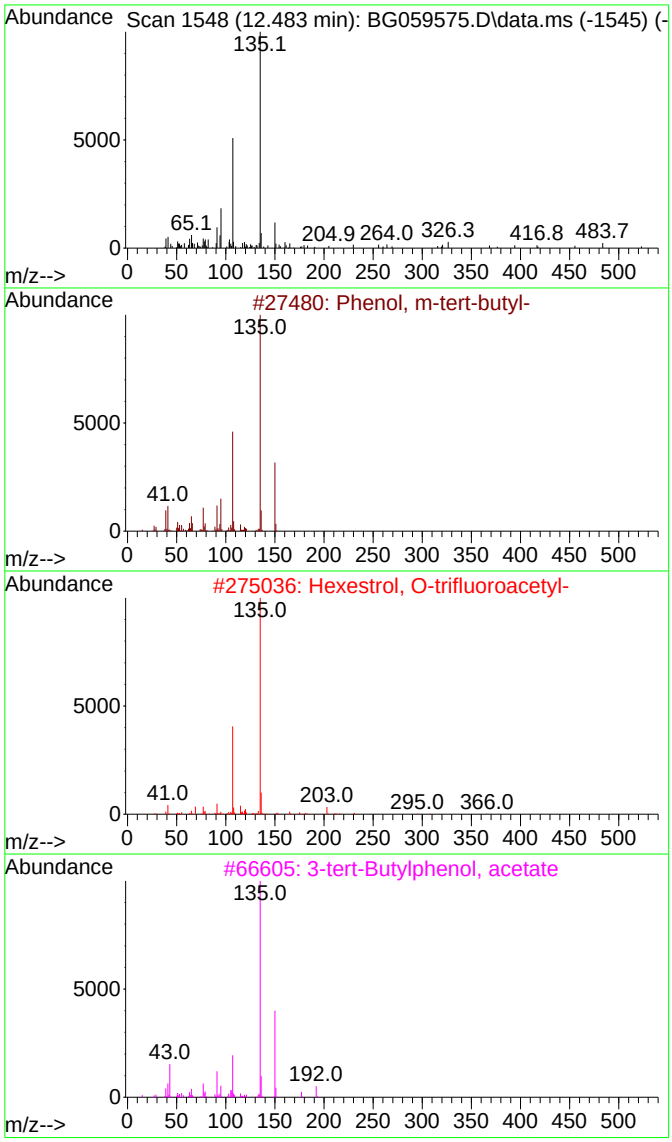
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

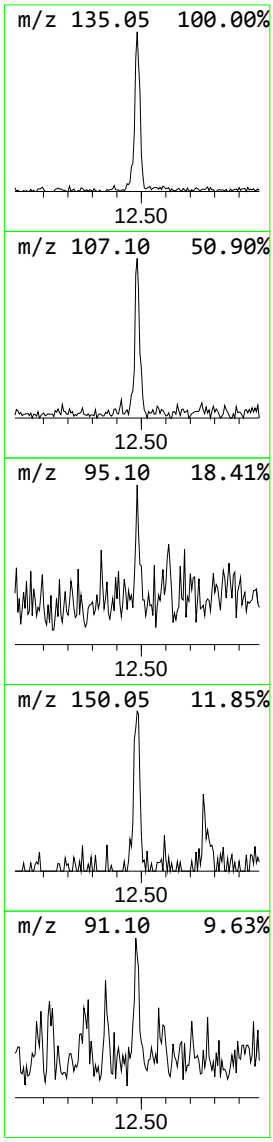
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Phenol, m-tert-butyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.483	2.99 ng/ul	114708	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	86
2	Hexestrol, O-trifluoroacetyl-	366	C20H21F3O3	1000365-31-7	64
3	3-tert-Butylphenol, acetate	192	C12H16O2	013189-51-0	59
4	4-Hydroxy-3-methylacetophenone	150	C9H10O2	000876-02-8	58
5	Cyclohexene, 2-ethenyl-1,3,3-tri...	150	C11H18	005293-90-3	53





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Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

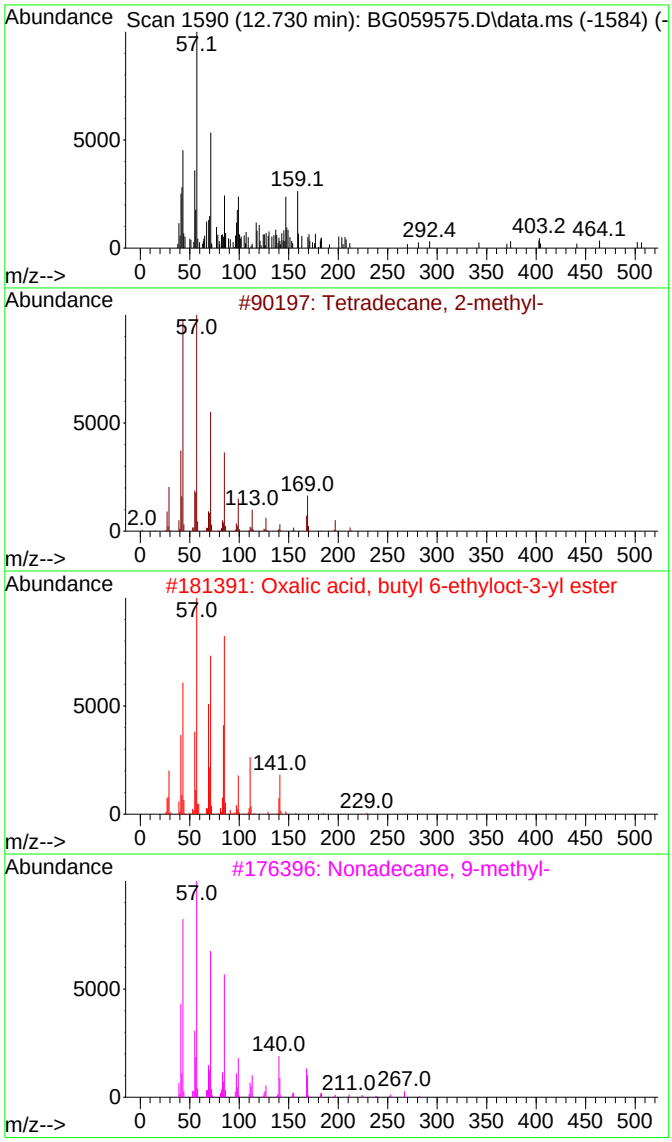
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

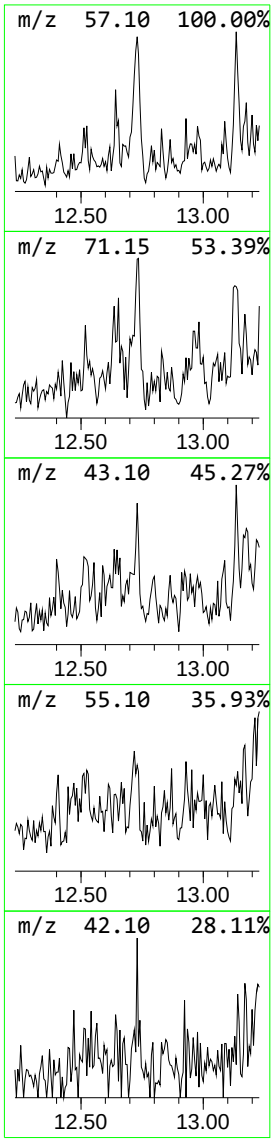
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.730	3.06 ng/ul	117579	Naphthalene-d8	11.208

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 2-methyl-	212	C15H32	001560-95-8	43
2	Oxalic acid, butyl 6-ethyloct-3-...	286	C16H30O4	1000309-34-2	38
3	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	35
4	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	35
5	Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	35





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

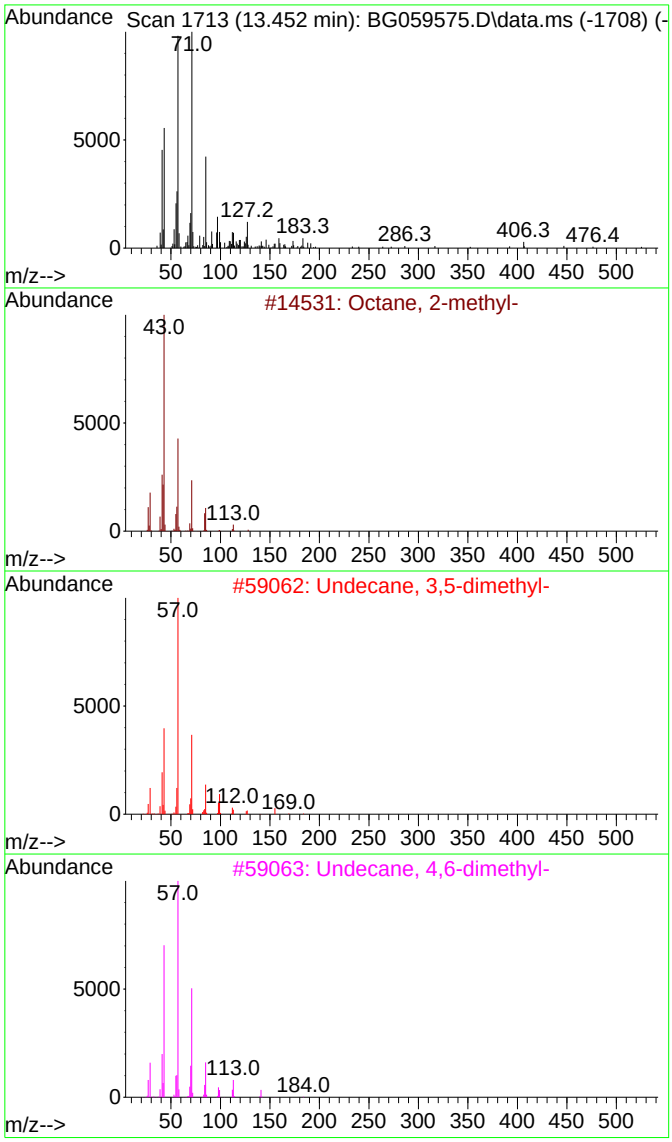
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

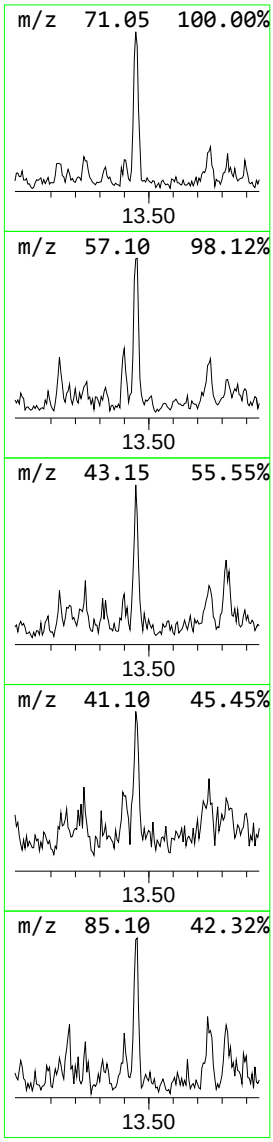
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.452	5.28 ng/ul	228032	Acenaphthene-d10	14.986

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	80
2	Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	64
3	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	62
4	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	59
5	Sulfurous acid, hexadecyl pentyl...	376	C21H44O3S	1000309-14-9	53





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

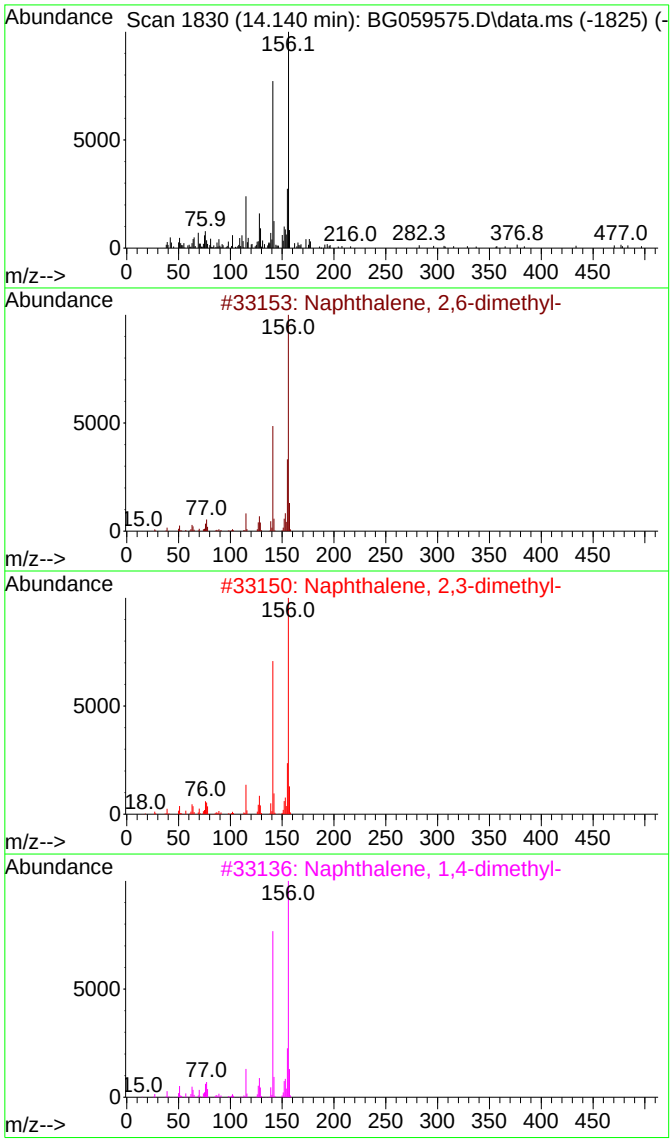
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

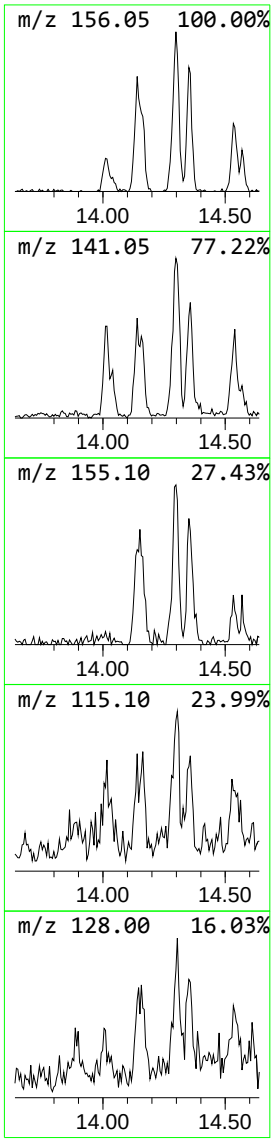
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Naphthalene, 2,6-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.140	3.08 ng/ul	132957	Acenaphthene-d10	14.986

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	93
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	93
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

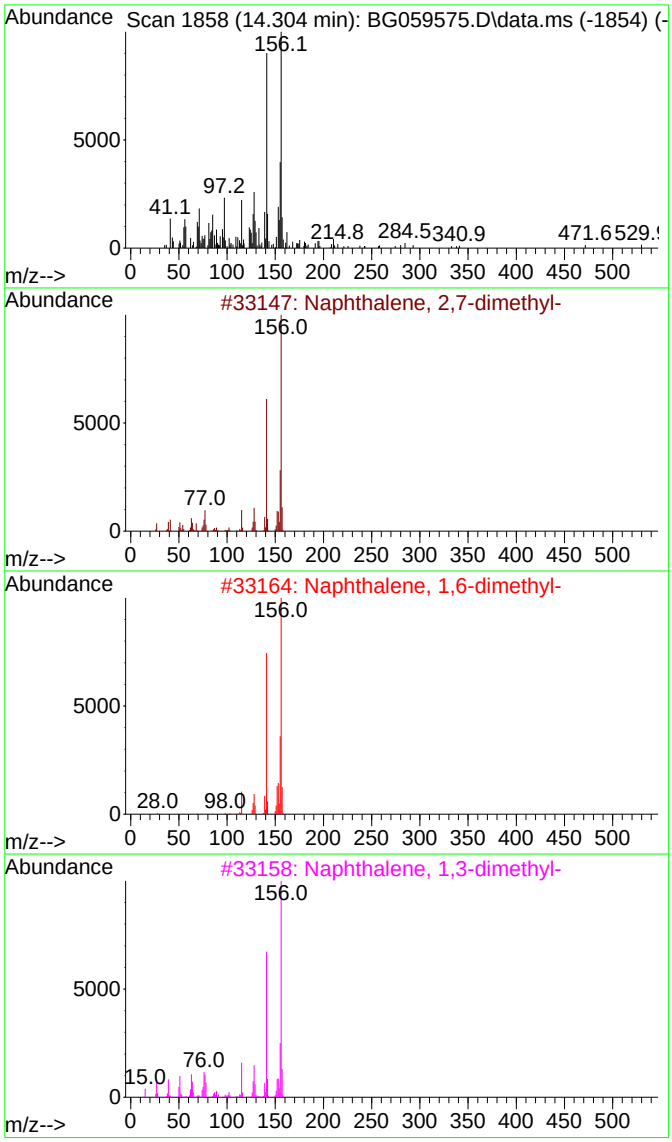
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

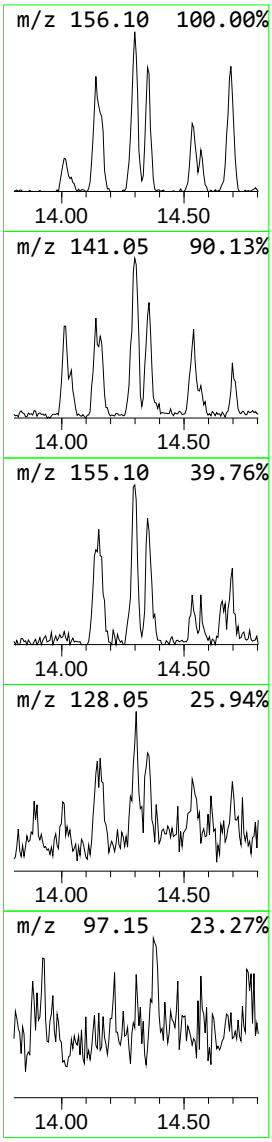
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Naphthalene, 2,7-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.304	5.05 ng/ul	217894	Acenaphthene-d10	14.986

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	87
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	83
3	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	81
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	81
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	81





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

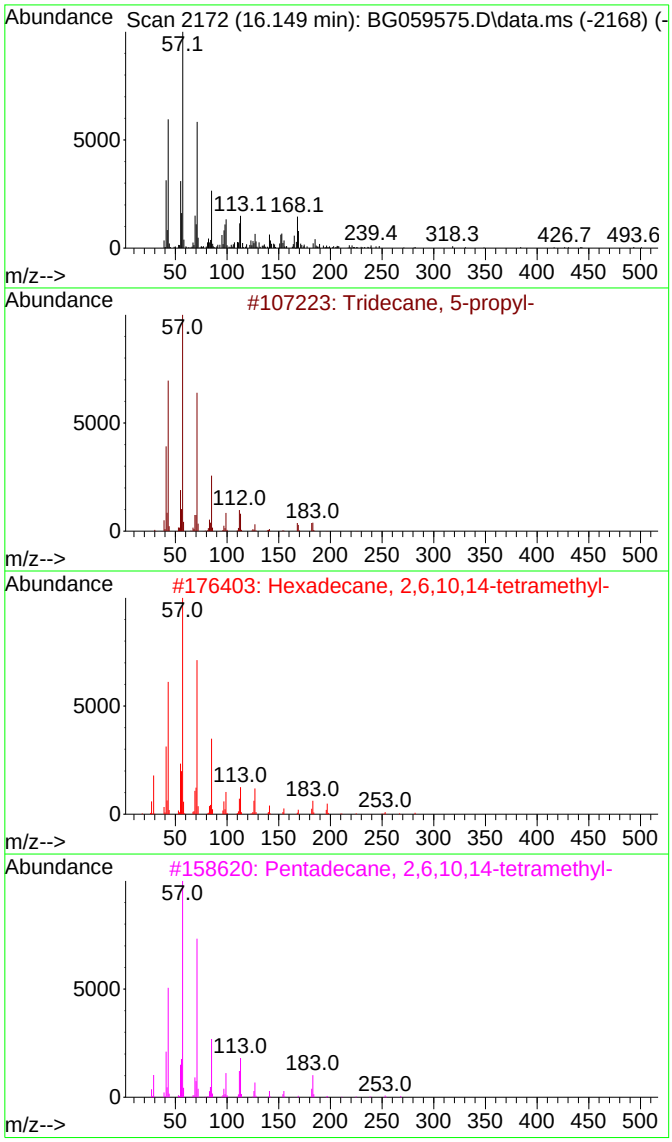
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

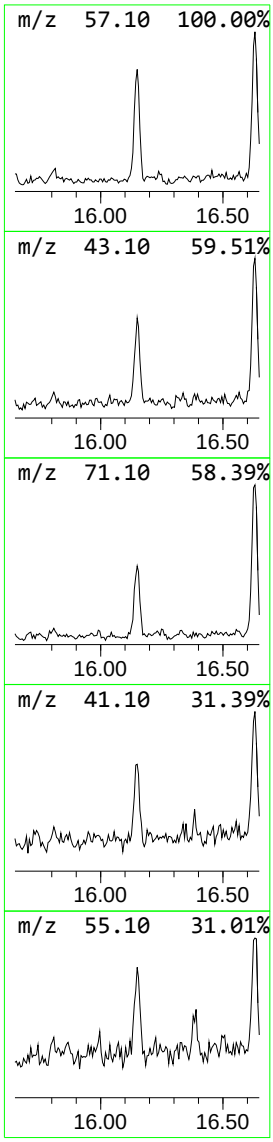
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.149	12.65 ng/ul	546348	Acenaphthene-d10	14.986

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-	226	C16H34	055045-11-9	91
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	76
3	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	72
4	Decane, 2-methyl-	156	C11H24	006975-98-0	68
5	Hexadecane	226	C16H34	000544-76-3	64





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

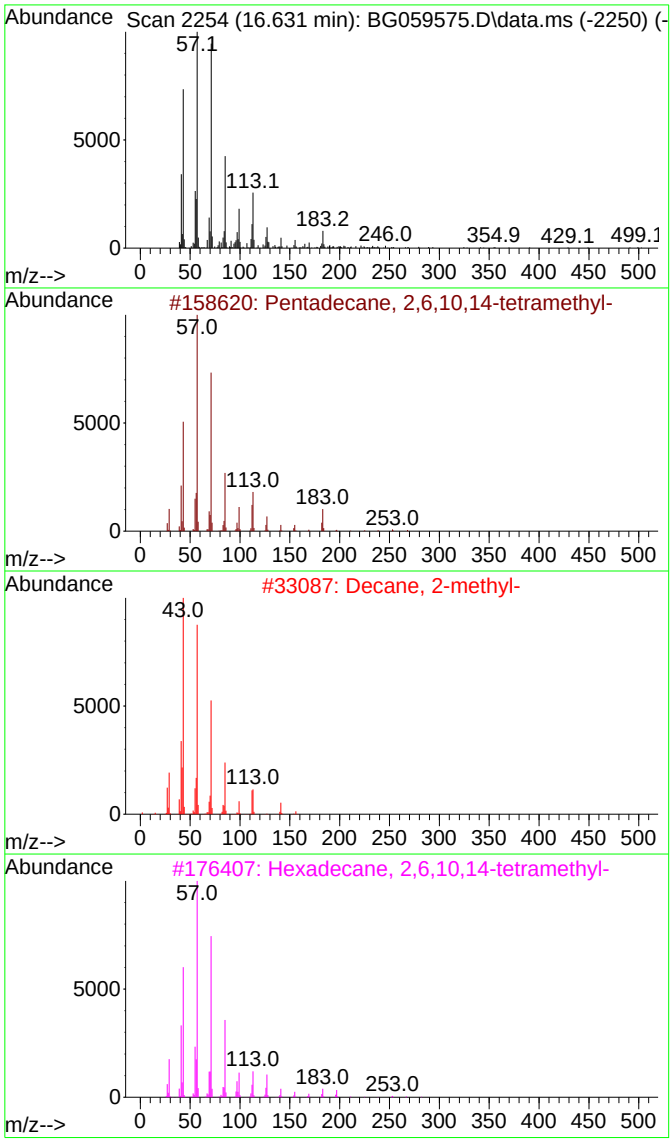
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

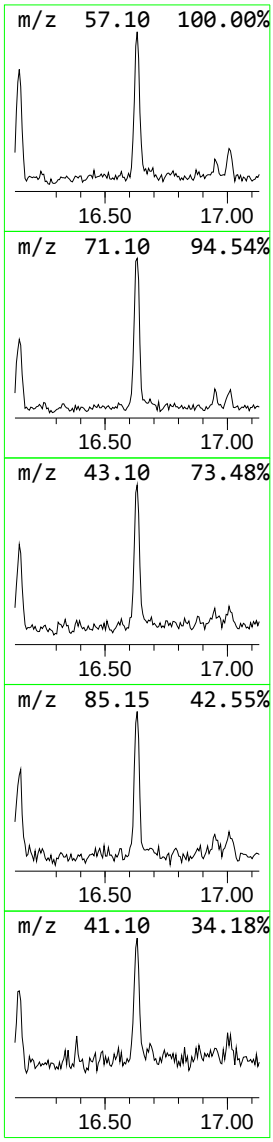
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.631	15.56 ng/ul	650332	Phenanthrene-d10	17.735

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
2	Decane, 2-methyl-	156	C11H24	006975-98-0	86
3	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	86
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	83
5	Heptadecane, 4-methyl-	254	C18H38	026429-11-8	81





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Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

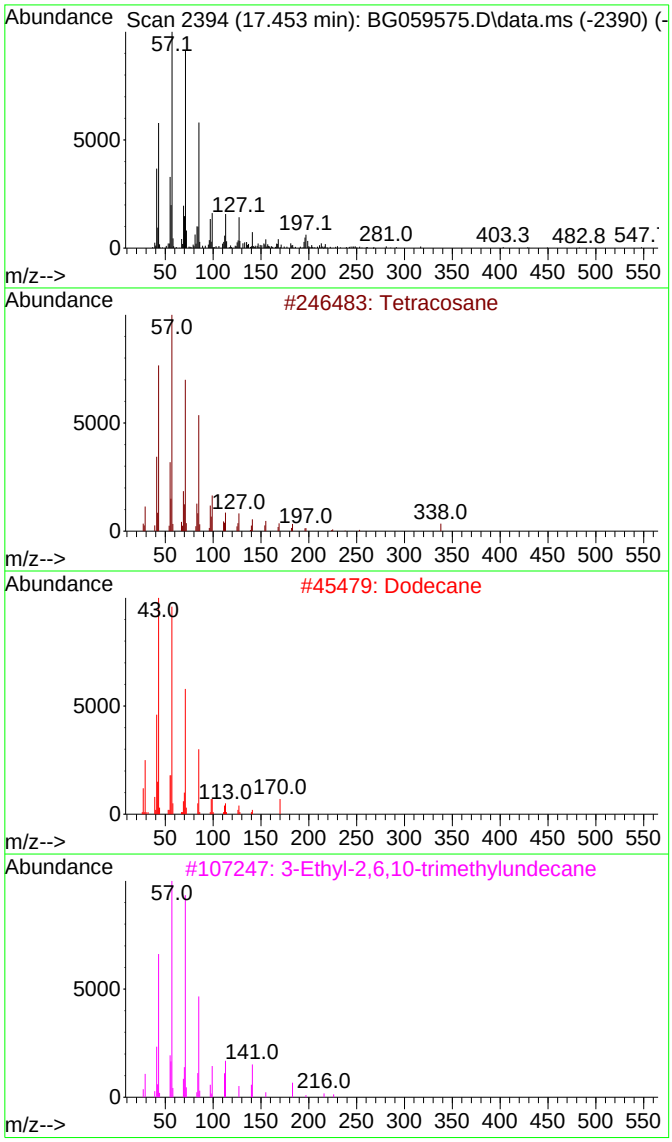
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
Quant Title : SVOA CALIBRATION

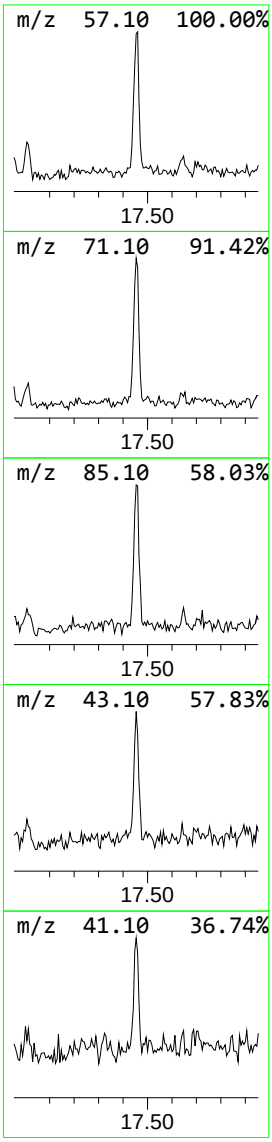
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.453	14.36 ng/ul	600494	Phenanthrene-d10	17.735

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	90
2	Dodecane	170	C12H26	000112-40-3	87
3	3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	86
4	2-Bromotetradecane	276	C14H29Br	074036-95-6	86
5	Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110123\
Data File : BG059575.D
Acq On : 1 Nov 2023 11:57
Operator : MA/JU
Sample : 04922-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EOAQ3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.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: C...	6.925	2.6	ng/ul	70252	1	8.358	544189	20.0
unknown-01	7.295	3.0	ng/ul	80173	1	8.358	544189	20.0
3-Bromooctane	8.247	4.0	ng/ul	108819	1	8.358	544189	20.0
Benzene, (1-eth...	8.717	4.5	ng/ul	122122	1	8.358	544189	20.0
Benzene, 1,4-di...	8.816	3.0	ng/ul	81131	1	8.358	544189	20.0
unknown-02	8.875	2.7	ng/ul	73088	1	8.358	544189	20.0
Bicyclo[4.2.0]o...	8.981	4.3	ng/ul	118061	1	8.358	544189	20.0
Benzene, 2-ethy...	9.439	6.7	ng/ul	181227	1	8.358	544189	20.0
unknown-03	9.686	2.8	ng/ul	76197	1	8.358	544189	20.0
unknown-04	9.774	2.3	ng/ul	62626	1	8.358	544189	20.0
Benzene, 1,2,3,...	9.974	2.5	ng/ul	97470	2	11.208	767749	20.0
unknown-05	10.350	5.4	ng/ul	206779	2	11.208	767749	20.0
unknown-06	10.562	6.3	ng/ul	240019	2	11.208	767749	20.0
(DEL) Alkane: S...	11.267	6.0	ng/ul	231450	2	11.208	767749	20.0
2-Ethyl-2,3-dih...	11.836	6.2	ng/ul	237900	2	11.208	767749	20.0
unknown-07	11.936	2.1	ng/ul	80593	2	11.208	767749	20.0
(DEL) Alkane: S...	12.118	4.9	ng/ul	186390	2	11.208	767749	20.0
Phenol, m-tert-...	12.483	3.0	ng/ul	114708	2	11.208	767749	20.0
(DEL) Alkane: S...	12.730	3.1	ng/ul	117579	2	11.208	767749	20.0
(DEL) Alkane: S...	13.452	5.3	ng/ul	228032	3	14.986	863547	20.0
Naphthalene, 2,...	14.140	3.1	ng/ul	132957	3	14.986	863547	20.0
Naphthalene, 2,...	14.304	5.0	ng/ul	217894	3	14.986	863547	20.0
(DEL) Alkane: S...	16.149	12.7	ng/ul	546348	3	14.986	863547	20.0
(DEL) Alkane: S...	16.631	15.6	ng/ul	650332	4	17.735	836113	20.0
(DEL) Alkane: S...	17.453	14.4	ng/ul	600494	4	17.735	836113	20.0