LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : GB863

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-BG110321.M
Title : SVOA CALIBRATION

Signal : TIC: BG050911.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.585	12	16	22	rVV	11617	17198	0.89%	0.045%
2	3.937	70	76	84	rBV4	4401	8171	0.42%	0.021%
3	4.014	84	89	109	rVB	145816	269490	13.90%	0.703%
4	4.255	124	130	134	rBV2	3318	5638	0.29%	0.015%
5	4.355	143	147	153	rBV4	4615	8407	0.43%	0.022%
6	4.625	184	193	197	rBV5	9386	19755	1.02%	0.052%
7	4.684	201	203	210	rVB4	6339	9941	0.51%	0.026%
8	5.295	302	307	315	rVB	33730	51602	2.66%	0.135%
9	5.512	338	344	350	rBV2	23167	38095	1.97%	0.099%
10	6.388	486	493	502	rVV	10190	16746	0.86%	0.044%
11	7.069	601	609	611	rBV5	3142	5481	0.28%	0.014%
12	7.369	651	660	670	rBV2	282376	544977	28.11%	1.422%
13	7.545	683	690	697	rBV	143257	250216	12.91%	0.653%
14	7.757	719	726	729	rBV	182088	341324	17.61%	0.891%
15	7.792	729	732	741	rVB	221542	359689	18.55%	0.938%
16	8.232	800	807	817	rVB	121310	209919	10.83%	0.548%
17	8.755	886	896	902	rBV	241427	606830	31.30%	1.583%
18	8.802	902	904	912	rVB	24130	35917	1.85%	0.094%
19	8.932	918	926	938	rVB	217055	394475	20.35%	1.029%
20	9.055	938	947	957	rBV	296262	533358	27.51%	1.392%
21	9.325	986	993	999	rBV	96358	172288	8.89%	0.450%
22	9.408	999	1007	1010	rVV	174212	344294	17.76%	0.898%
23	9.449	1010	1014	1023	rVB	333401	574579	29.64%	1.499%
24	9.701	1047	1057	1064	rBV	421963	705748	36.40%	1.841%
25	10.130	1123	1130	1132	rBV	156258	267150	13.78%	0.697%
26	10.160	1132	1135	1150	rVB	229419	435023	22.44%	1.135%
27	10.671	1214	1222	1236	rBV2	228271	681527	35.16%	1.778%
28	10.941	1262	1268	1274	rBV2	12437	23957	1.24%	0.063%
29	11.059	1280	1288	1292	rBV	170724	314080	16.20%	0.819%
30	11.106	1292	1296	1303	rVV	157900	284934	14.70%	0.743%
31	11.188	1303	1310	1321	rVB	99396	197986	10.21%	0.517%
32	11.746	1399	1405	1413	rVB	11067	16777	0.87%	0.044%
33	12.310	1495	1501	1510	rBV	379291	658920	33.99%	1.719%
34	12.551	1536	1542	1551	rBV	253158	435956	22.49%	1.137%
35	12.621	1553	1554	1561	rVB3	3527	4404	0.23%	0.011%
36	12.827	1584	1589	1594	rBV2	2883	4847	0.25%	0.013%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

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-BG110321.M
Title : SVOA CALIBRATION

37	12.915	1596	1604	1612	rVV	540109	914149	47.15%	2.385%
38	13.056	1621	1628	1636	rVB	186522	305907	15.78%	0.798%
39	13.291	1661	1668	1675	rBV	244898	421422	21.74%	1.100%
40	13.691	1729	1736	1743	rBV	565171	893403	46.08%	2.331%
41	13.943	1771	1779	1790	rBV	419495	699260	36.07%	1.824%
42	14.032	1790	1794	1799	rVB	10827	16021	0.83%	0.042%
43	14.249	1824	1831	1841	rBV	402909	584872	30.17%	1.526%
44	14.337	1841	1846	1854	rVB	89996	133676	6.90%	0.349%
45	14.425	1854	1861	1869	rVB	347216	519569	26.80%	1.356%
46	14.554	1876	1883	1885	rVV	469452	764350	39.43%	1.994%
47	14.584	1885	1888	1895	rVB	660908	992724	51.21%	2.590%
48	14.854	1924	1934	1940	rBV	268229	429958	22.18%	1.122%
49	14.919	1940	1945	1953	rVB	69082	108122	5.58%	0.282%
50	15.054	1959	1968	1988	rBV3	437442	957797	49.41%	2.499%
51	15.212	1988	1995	1999	rVV	416241	679002	35.02%	1.772%
52	15.254	1999	2002	2010	rVB	274603	394099	20.33%	1.028%
53	15.495	2036	2043	2050	rVB	95322	149748	7.72%	0.391%
54	15.653	2063	2070	2077	rBV	672510	940903	48.53%	2.455%
55	15.847	2095	2103	2108	rBV	550668	873558	45.06%	2.279%
56	15.900	2108	2112	2117	rVV	133468	202769	10.46%	0.529%
57	15.959	2117	2122	2135	rVB	187392	275632	14.22%	0.719%
58	16.100	2135	2146	2153	rBV	856504	1313488	67.75%	3.427%
59	16.335	2181	2186	2193	rBV3	17750	33909	1.75%	0.088%
60	16.470	2203	2209	2213	rVB	10314	14485	0.75%	0.038%
61	16.687	2243	2246	2250	rVV3	3548	5398	0.28%	0.014%
62	16.769	2250	2260	2264	rVV5	10632	20326	1.05%	0.053%
63	16.840	2264	2272	2277	rVV2	163656	254210	13.11%	0.663%
64	16.899	2277	2282	2288	rVB	457060	698491	36.03%	1.822%
65	17.245	2336	2341	2355	rBV2	245241	422707	21.80%	1.103%
66	17.604	2395	2402	2405	rVV	340991	559269	28.85%	1.459%
67	17.645	2405	2409	2414	rVV	620853	943008	48.64%	2.460%
68	17.698	2414	2418	2425	rVB	615651	928612	47.90%	2.423%
69	18.203	2499	2504	2509	rBV	36629	48660	2.51%	0.127%
70	18.538	2553	2561	2570	rBV	780749	1034827	53.38%	2.700%
71	18.826	2606	2610	2616	rVB3	7881	9784	0.50%	0.026%
72	19.543	2728	2732	2738	rVB4	7741	11888	0.61%	0.031%
73	19.613	2738	2744	2750	rBV	8506	12425	0.64%	0.032%
74	19.836	2776	2782	2788	rBV	79716	108692	5.61%	0.284%
75	19.883	2788	2790	2796	rVB2	4089	4874	0.25%	0.013%
76	19.977	2799	2806	2809	rBV2	744476	1266419	65.33%	3.304%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

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-BG110321.M
Title : SVOA CALIBRATION

77	20.189	2837	2842	2848	rBV	95805	125950	6.50%	0.329%
78	20.342	2863	2868	2872	rVV6	5449	7718	0.40%	0.020%
79	20.806	2943	2947	2950	rBV4	3494	6076	0.31%	0.016%
80	20.994	2974	2979	2984	rVB	57700	75675	3.90%	0.197%
81	21.117	2996	3000	3006	rBV5	3128	5644	0.29%	0.015%
82	21.164	3006	3008	3014	rBV6	3920	4346	0.22%	0.011%
83	21.787	3104	3114	3122	rVB2	29787	52398	2.70%	0.137%
84	21.881	3122	3130	3141	rBV2	888719	1938634	100.00%	5.058%
85	22.063	3154	3161	3164	rVB7	3492	5491	0.28%	0.014%
86	22.245	3186	3192	3202	rVB2	44284	77818	4.01%	0.203%
87	22.457	3221	3228	3237	rBV	834691	1465302	75.58%	3.823%
88	22.645	3254	3260	3265	rBV4	43553	76911	3.97%	0.201%
89	23.021	3316	3324	3334	rVB	484190	889655	45.89%	2.321%
90	23.421	3386	3392	3400	rBV5	9710	22625	1.17%	0.059%
91	24.214	3516	3527	3535	rBV	425326	1093146	56.39%	2.852%
92	25.066	3660	3672	3679	rVV2	348339	1053536	54.34%	2.749%
93	25.142	3679	3685	3698	rVV2	228218	620466	32.01%	1.619%
94	25.301	3700	3712	3726	rVB2	186709	585369	30.19%	1.527%
95	26.476	3903	3912	3923	rBV3	23203	67446	3.48%	0.176%
96	27.921	4148	4158	4167	rVV6	21276	75525	3.90%	0.197%
97	29.278	4370	4389	4409	rBV2	319460	1510313	77.91%	3.940%
98	29.643	4446	4451	4452	rBV5	9412	12604	0.65%	0.033%
99	30.436	4565	4586	4605	rBV4	141906	739352	38.14%	1.929%
100	31.746	4801	4809	4810	rBV6	8878	18285	0.94%	0.048%

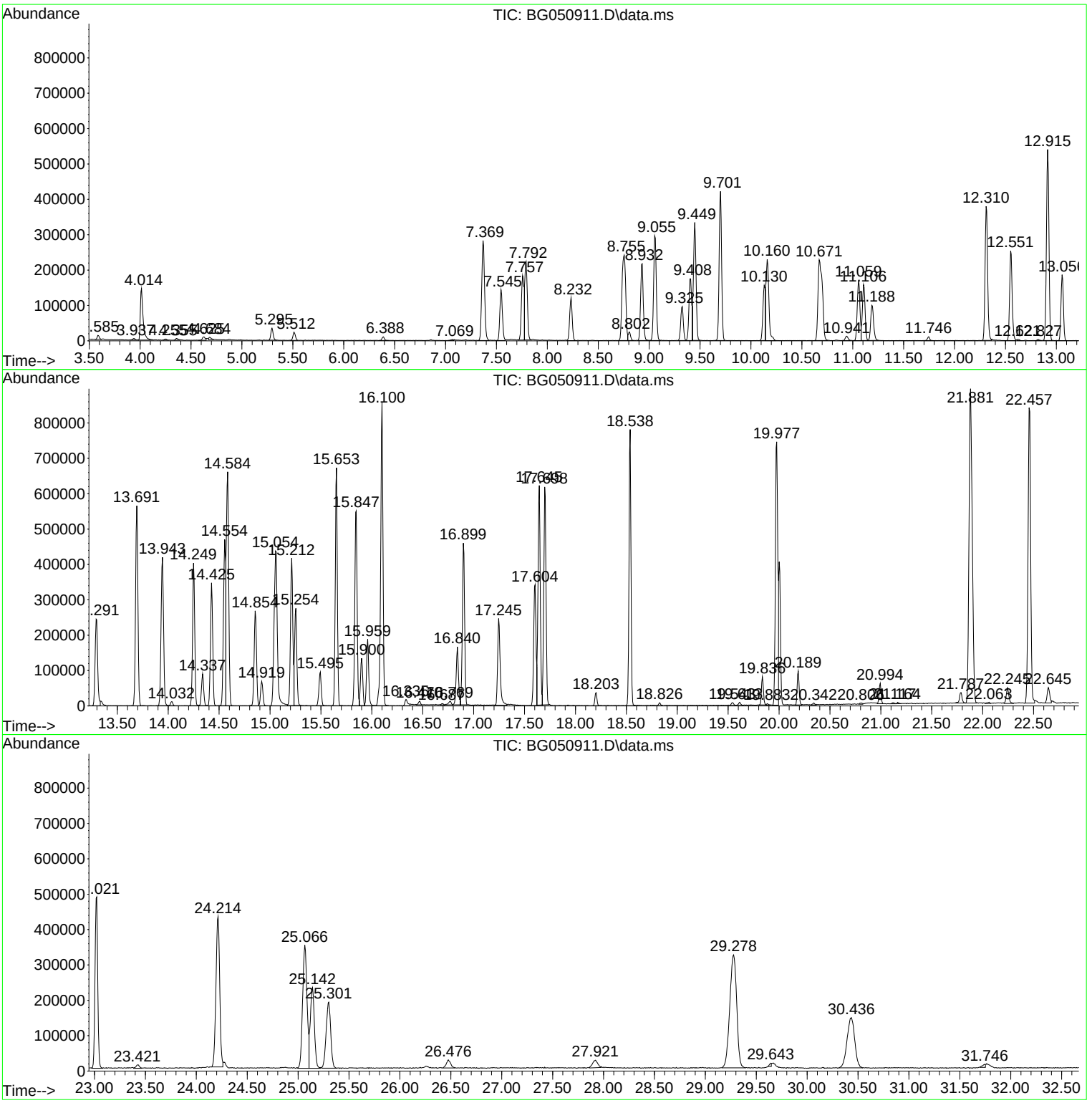
Sum of corrected areas: 38328397

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

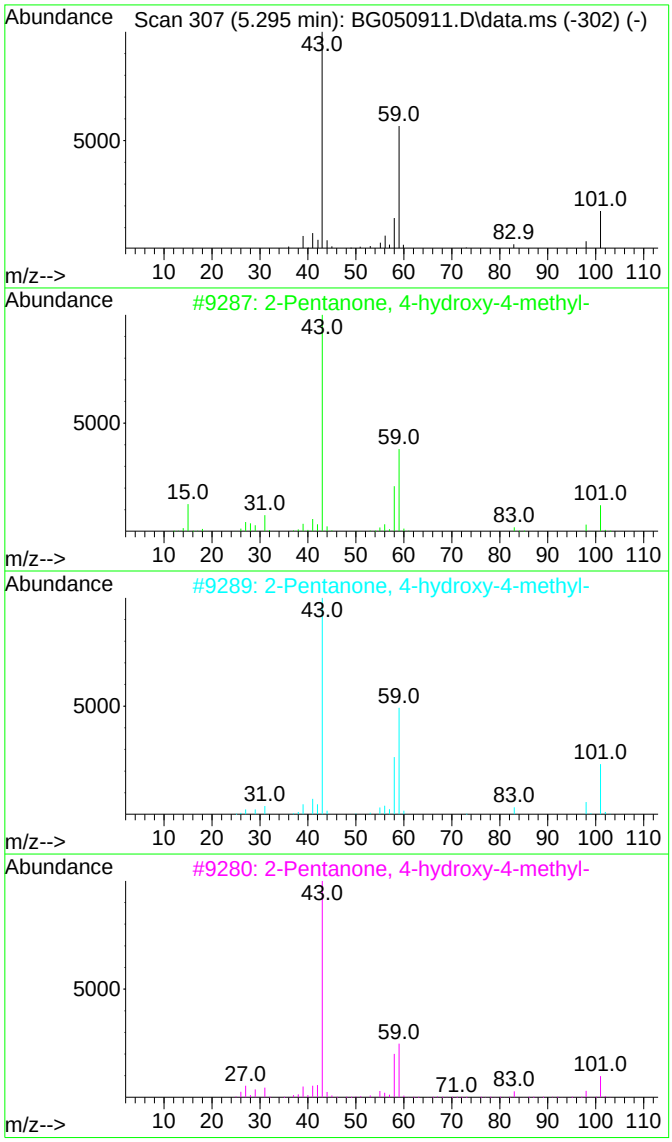
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

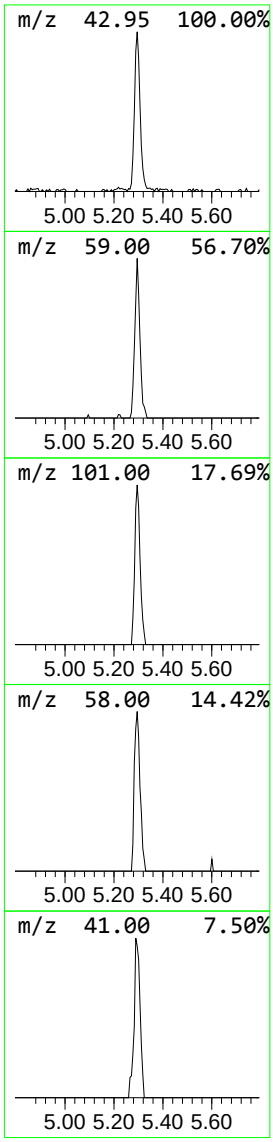
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.295	4.92 ng/ul	51602	1,4-Dichlorobenzene-d4	8.232

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
5	3-Octanol	130	C8H18O	000589-98-0	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

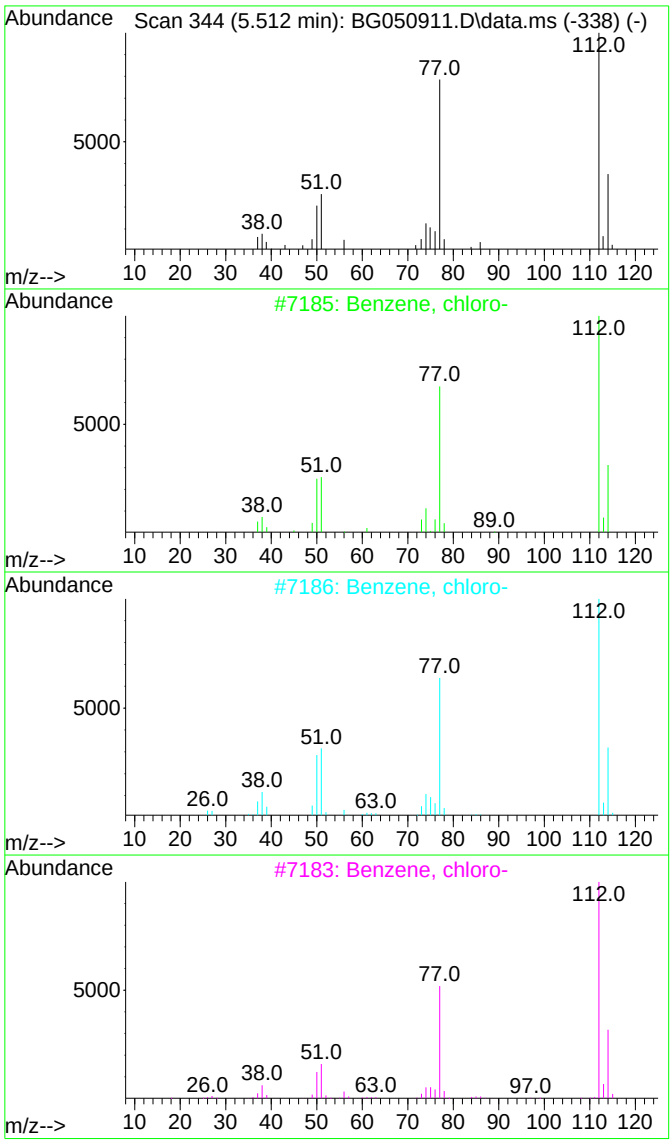
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

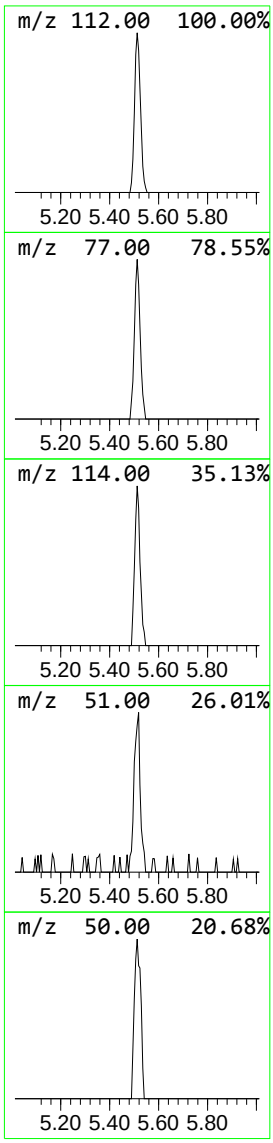
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.512	3.63 ng/ul	38095	1,4-Dichlorobenzene-d4	8.232

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

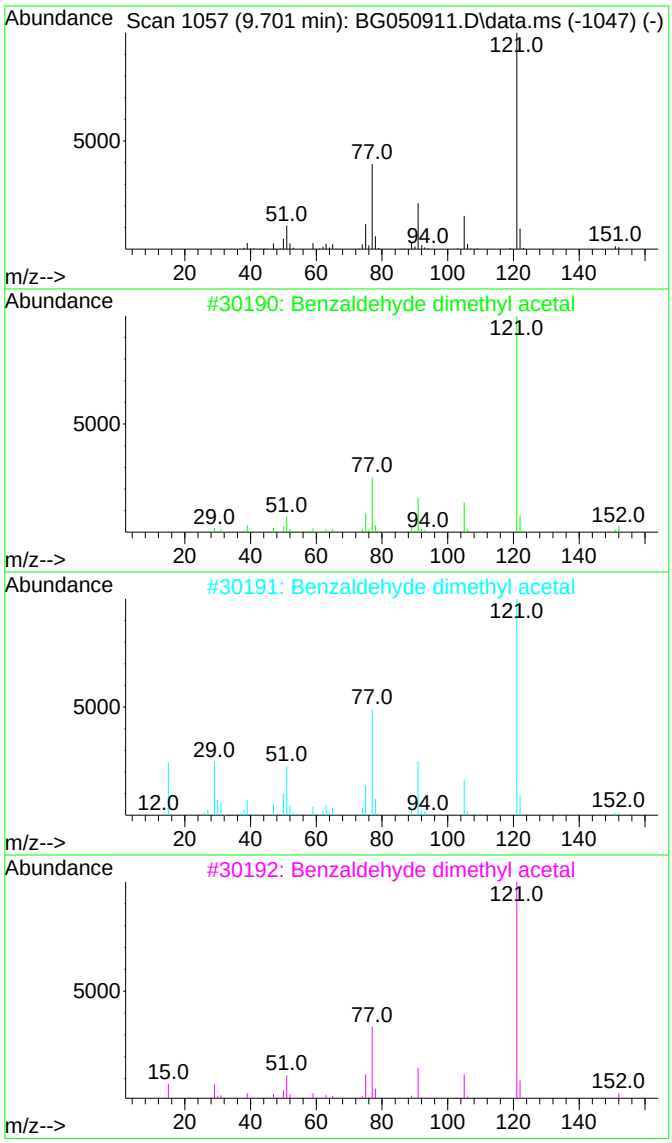
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

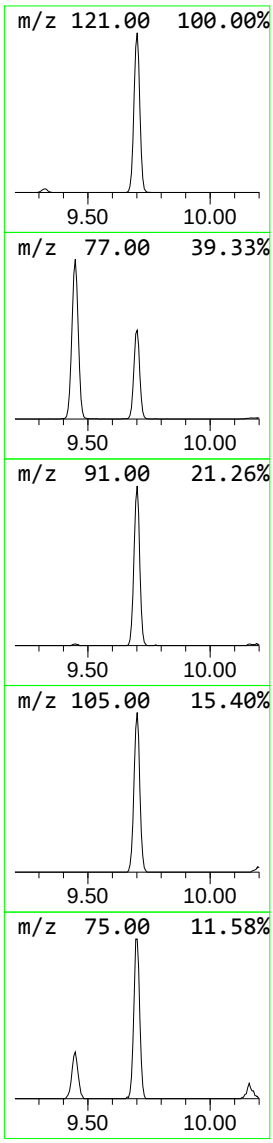
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzaldehyde dimethyl acetal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.701	44.94 ng/ul	705748	Naphthalene-d8	11.059

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	94
2	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	91
3	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	91
4	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	87
5	Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

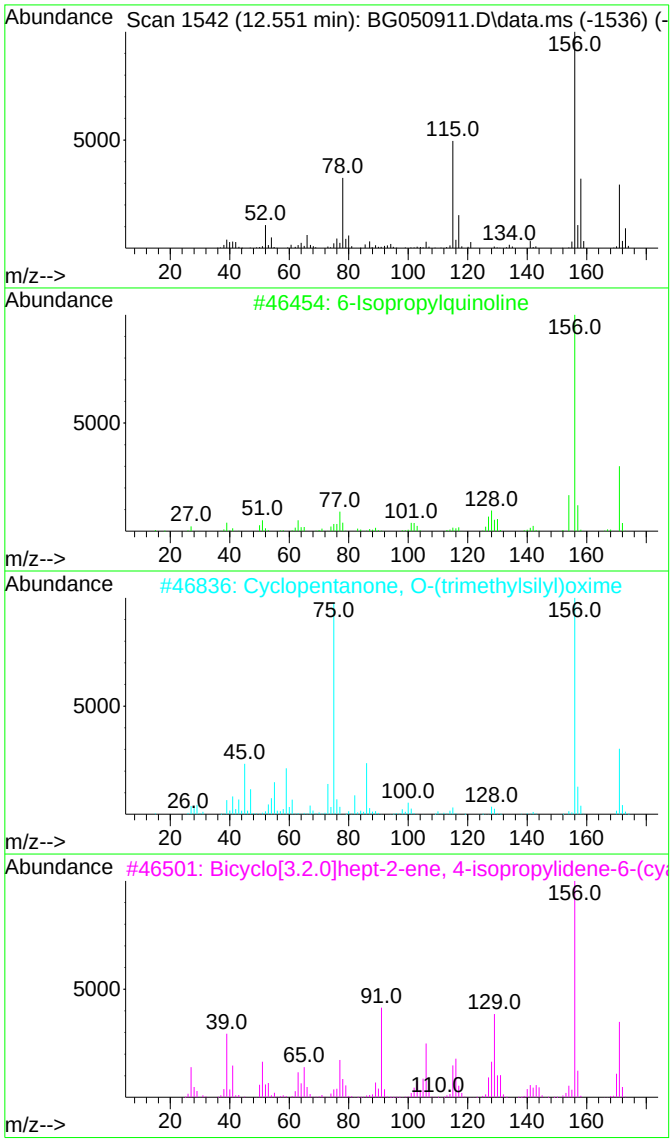
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

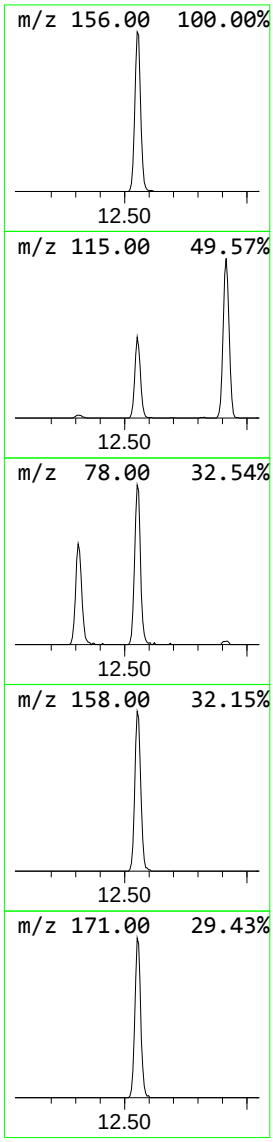
****

Peak Number 4 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.551	27.76 ng/ul	435956	Naphthalene-d8	11.059

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Isopropylquinoline	171	C12H13N	000135-79-5	38
2	Cyclopentanone, O-(trimethylsily...	171	C8H17NOSi	026208-87-7	32
3	Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32
4	Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28
5	Tebuthiuron	228	C9H16N4OS	034014-18-1	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

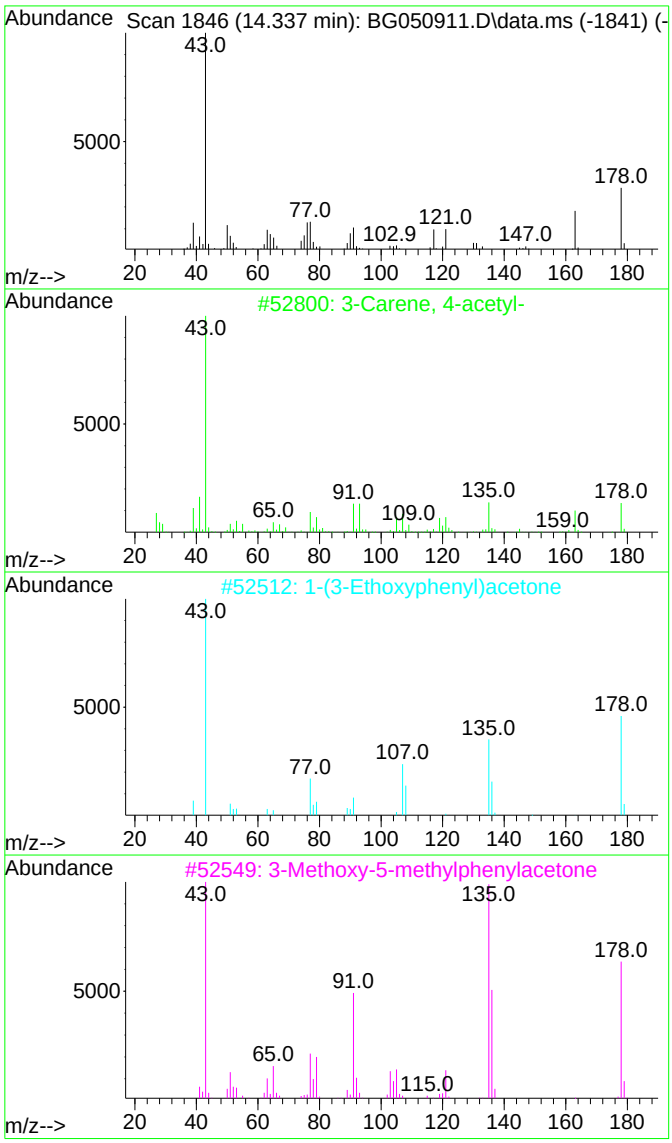
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

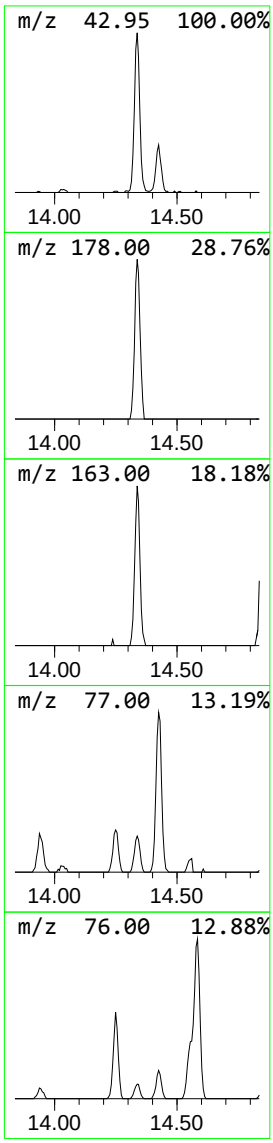
****

Peak Number 5 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.337	6.22 ng/ul	133676	Acenaphthene-d10	14.854

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Carene, 4-acetyl-	178	C12H18O	1000156-14-1	38
2	1-(3-Ethoxyphenyl)acetone	178	C11H14O2	023037-47-0	38
3	3-Methoxy-5-methylphenylacetone	178	C11H14O2	1000386-47-1	27
4	(+)-2-Carene, 2-acetyl-	178	C12H18O	1000151-09-7	17
5	Acetamide, N-(2-phenylethyl)-	163	C10H13NO	000877-95-2	10





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

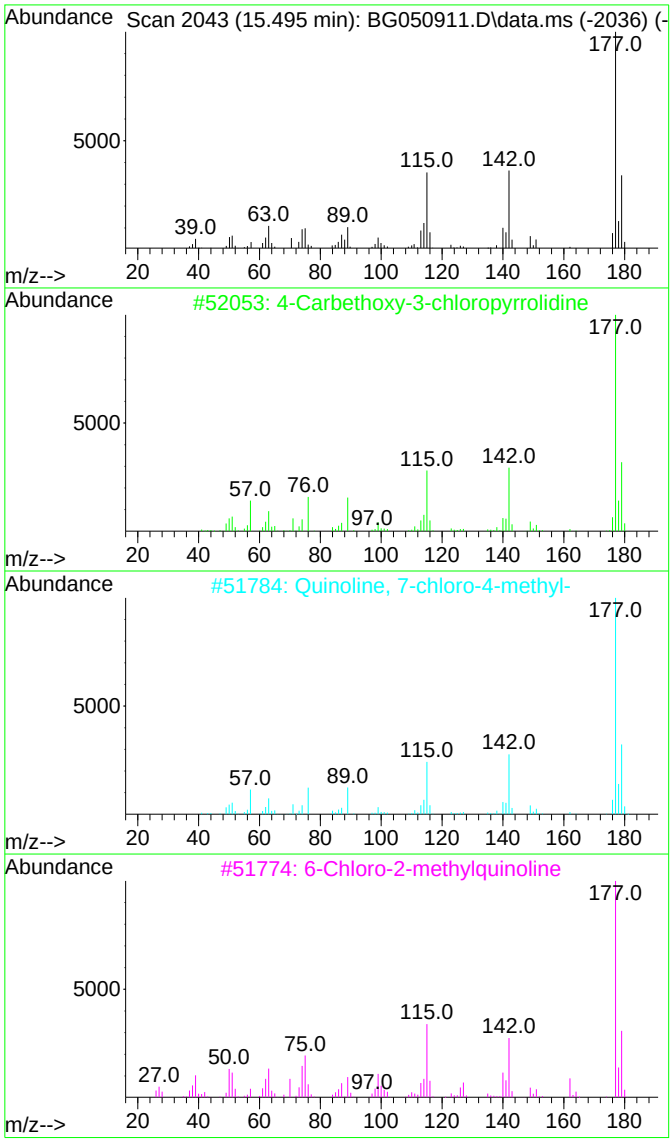
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

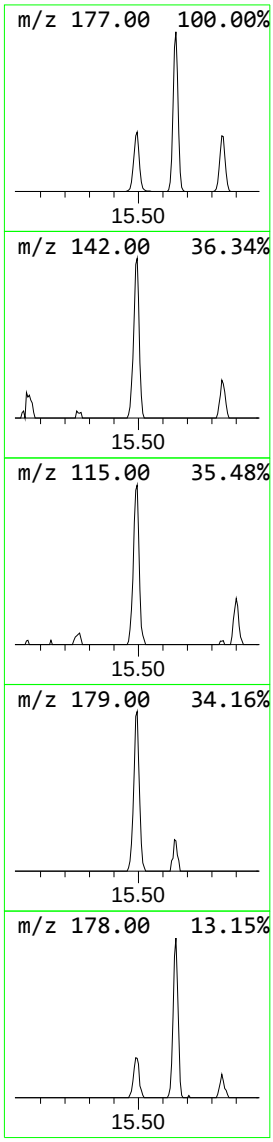
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4-Carbethoxy-3-chloropyrrol... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.495	6.97 ng/ul	149748	Acenaphthene-d10	14.854

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Carbethoxy-3-chloropyrrolidine	177	C7H12ClNO2	1010255-97-9	94
2	Quinoline, 7-chloro-4-methyl-	177	C10H8ClN	040941-53-5	94
3	6-Chloro-2-methylquinoline	177	C10H8ClN	000092-46-6	94
4	Quinoline, 2-chloro-4-methyl-	177	C10H8ClN	000634-47-9	91
5	2-Chloro-3-methylquinoline	177	C10H8ClN	057876-69-4	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

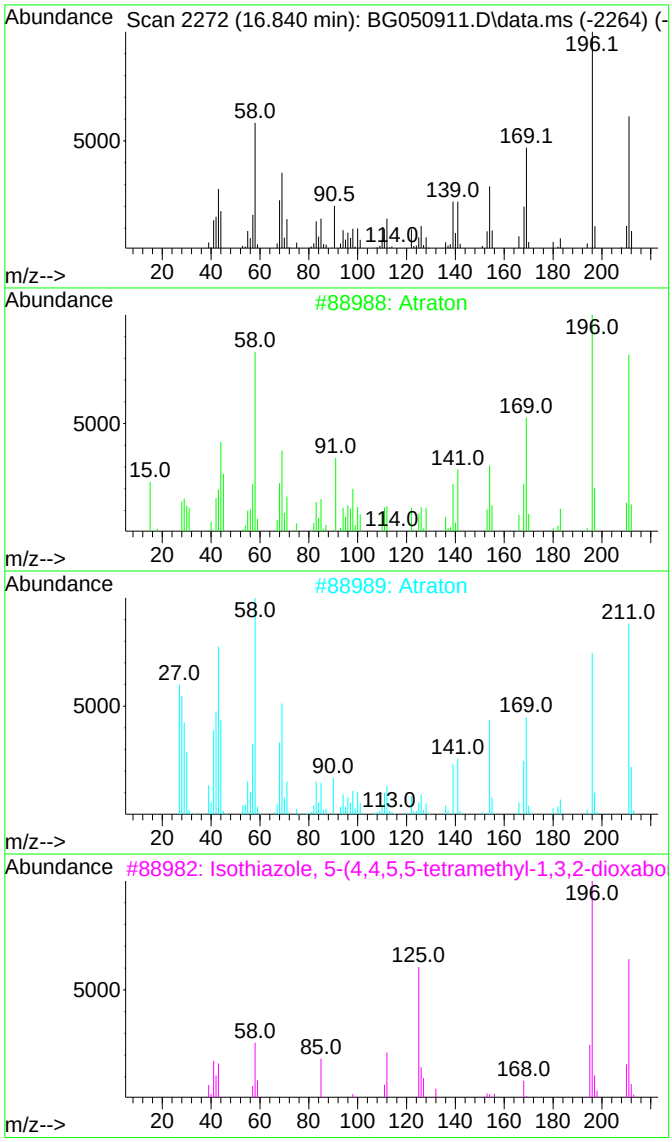
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

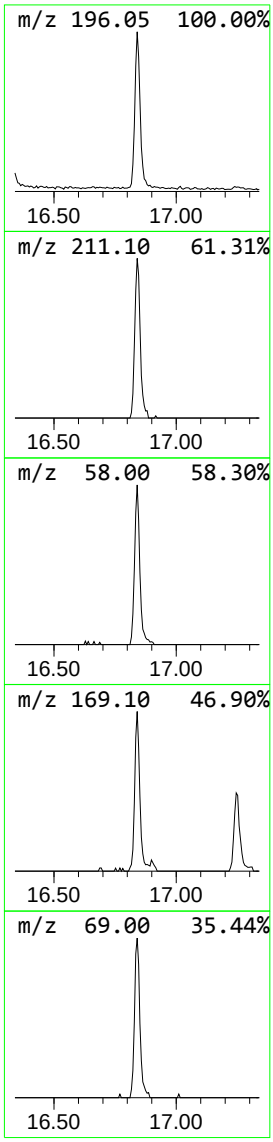
****

Peak Number 7 Atraton Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.840	9.09 ng/ul	254210	Phenanthrene-d10	17.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Atraton	211	C9H17N5O	001610-17-9	87
2	Atraton	211	C9H17N5O	001610-17-9	70
3	Isothiazole, 5-(4,4,5,5-tetramet...	211	C9H14BNO2S	1000337-80-0	41
4	N2,N2-Dimethyl-1,3,5-triazine-2,...	211	C8H17N5Si	1000471-09-5	38
5	2(1H)-Pyrimidinone,1,5-dimethyl-...	211	C9H17N3OSi	1000147-48-1	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

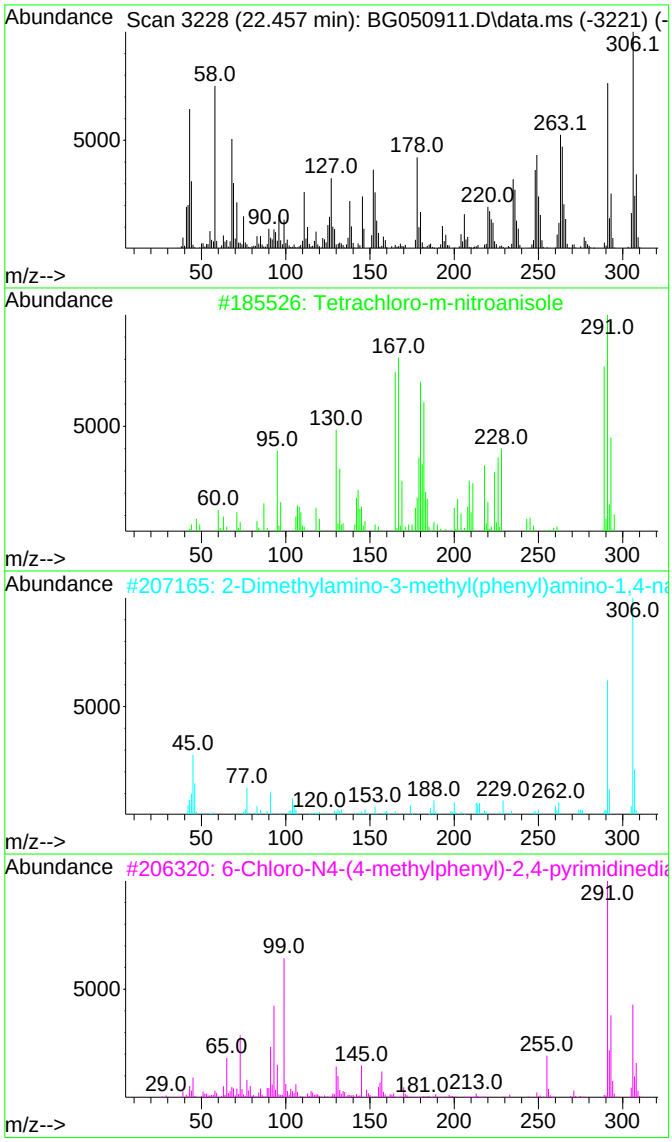
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

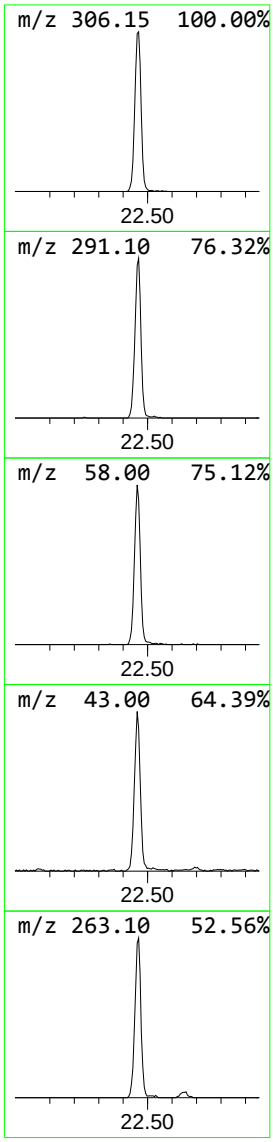
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Tetrachloro-m-nitroanisole Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.457	15.12 ng/ul	1465300	Chrysene-d12	21.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrachloro-m-nitroanisole	289	C7H3Cl4NO3	162627-34-1	59
2	2-Dimethylamino-3-methyl(phenyl)...	306	C19H18N2O2	134614-10-1	41
3	6-Chloro-N4-(4-methylphenyl)-2,4...	306	C14H19ClN4Si	1000486-00-3	27
4	1H-Carbazol-1-one, 2,3,4,9-tetra...	291	C19H17NO2	299405-79-1	25
5	1,1':2',1''-Terphenyl, 4'-phenyl-	306	C24H18	001165-53-3	22





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

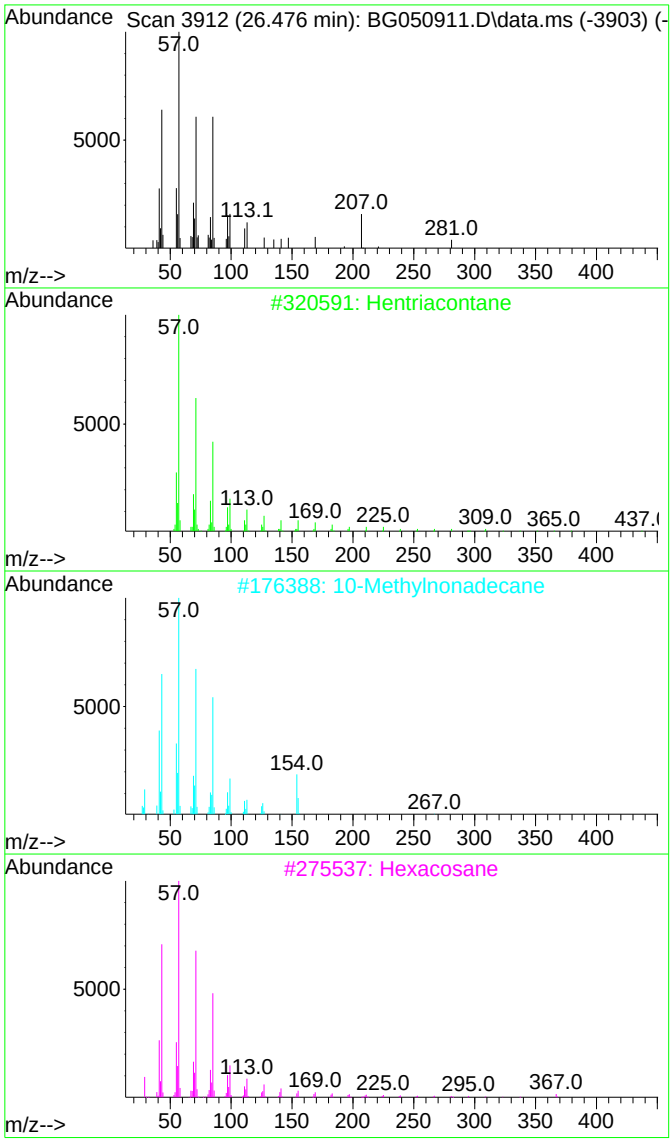
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

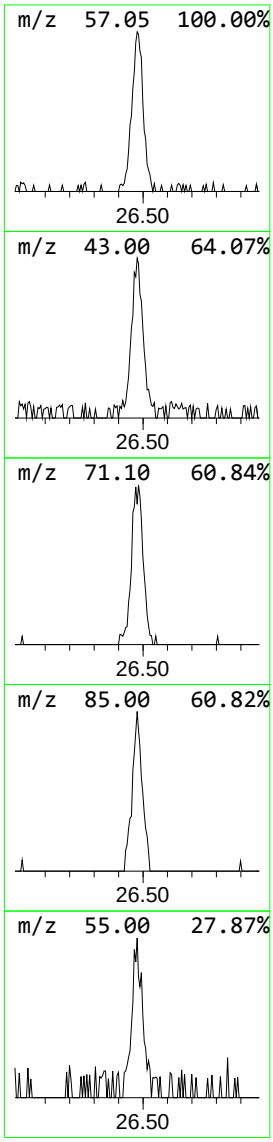
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.476	2.30 ng/ul	67446	Perylene-d12	25.301

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	86
2	10-Methylnonadecane	282	C20H42	056862-62-5	83
3	Hexacosane	366	C26H54	000630-01-3	80
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	64
5	Tetratetracontane	619	C44H90	007098-22-8	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

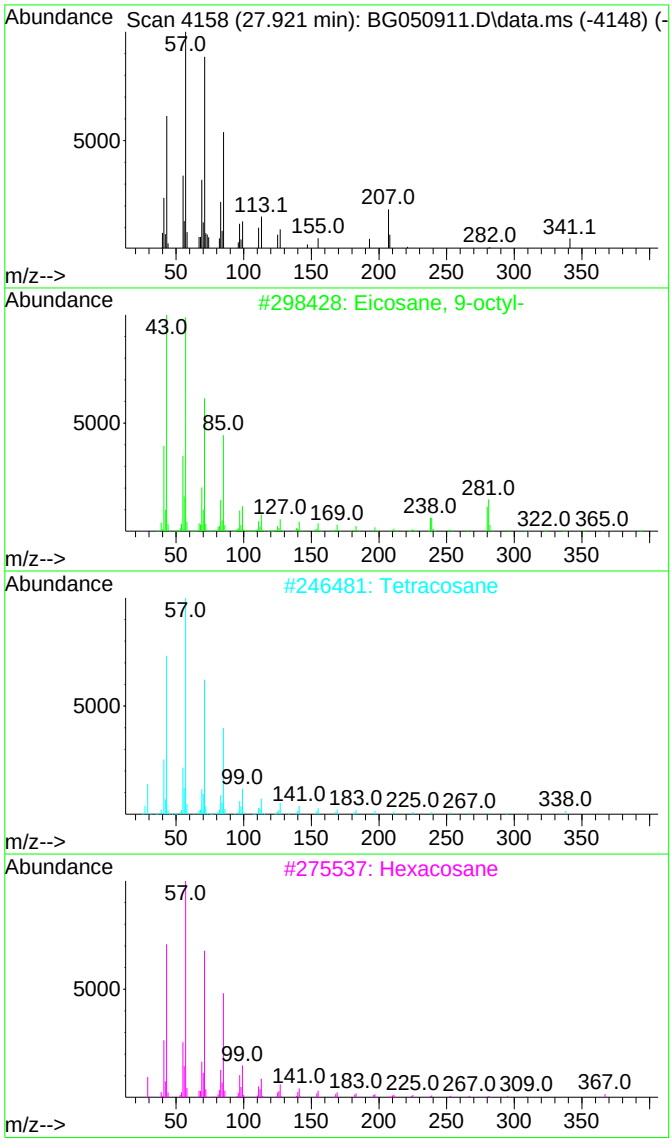
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

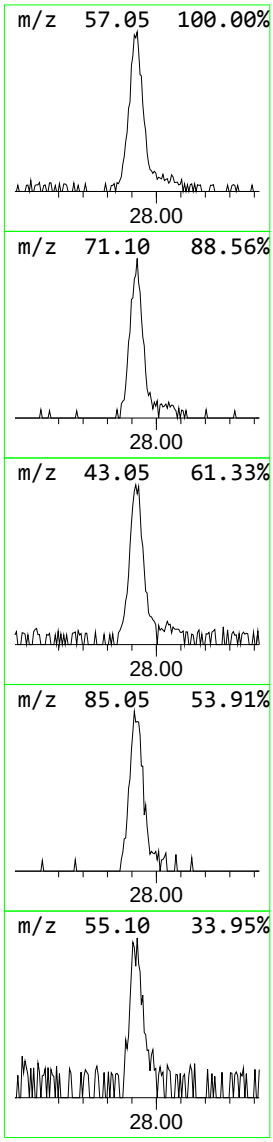
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.921	2.58 ng/ul	75525	Perylene-d12	25.301

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 9-octyl-	394	C28H58	013475-77-9	86
2	Tetracosane	338	C24H50	000646-31-1	80
3	Hexacosane	366	C26H54	000630-01-3	80
4	Pentacosane	352	C25H52	000629-99-2	72
5	Nonadecane	268	C19H40	000629-92-5	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
Data File : BG050911.D
Acq On : 9 Nov 2021 15:08
Operator : CG/JU
Sample : M4532-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GB863

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-...	5.295	4.9	ng/ul	51602	1	8.232	209919	20.0
Benzene, chloro-	5.512	3.6	ng/ul	38095	1	8.232	209919	20.0
Benzaldehyde di...	9.701	44.9	ng/ul	705748	2	11.059	314080	20.0
unknown-01	12.551	27.8	ng/ul	435956	2	11.059	314080	20.0
unknown-02	14.337	6.2	ng/ul	133676	3	14.854	429958	20.0
4-Carbethoxy-3-...	15.495	7.0	ng/ul	149748	3	14.854	429958	20.0
Atraton	16.840	9.1	ng/ul	254210	4	17.604	559269	20.0
Tetrachloro-m-n...	22.457	15.1	ng/ul	1465300	5	21.899	1938630	20.0
(DEL) Alkane: S...	26.476	2.3	ng/ul	67446	6	25.301	585369	20.0
(DEL) Alkane: S...	27.921	2.6	ng/ul	75525	6	25.301	585369	20.0