

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049648.D
 Acq On : 5 Aug 2021 13:53
 Operator : CG/JU
 Sample : M3162-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG07

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-BG080421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049648.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.073	3	6	16	rVB	46421	89127	3.55%	0.628%
2	3.156	16	20	29	rVB	45061	73212	2.92%	0.516%
3	3.426	62	66	72	rBV4	8297	12892	0.51%	0.091%
4	3.843	133	137	146	rBV	98584	185487	7.39%	1.308%
5	7.191	701	707	717	rVB	186424	344541	13.73%	2.429%
6	7.362	729	736	745	rVB	116553	203086	8.09%	1.432%
7	7.567	765	771	780	rBV	176060	310169	12.36%	2.187%
8	8.043	845	852	858	rBV	92552	166022	6.62%	1.171%
9	8.748	965	972	981	rBV	234715	412785	16.45%	2.910%
10	9.212	1042	1051	1063	rVB	185381	335730	13.38%	2.367%
11	9.612	1117	1119	1127	rVB3	3085	4262	0.17%	0.030%
12	9.935	1166	1174	1183	rBV	169873	319672	12.74%	2.254%
13	10.481	1259	1267	1276	rBV	230036	427108	17.02%	3.011%
14	10.622	1285	1291	1298	rBV	26890	50609	2.02%	0.357%
15	10.863	1323	1332	1340	rVB	132112	238791	9.51%	1.684%
16	10.998	1348	1355	1365	rBV	217812	404595	16.12%	2.853%
17	11.909	1503	1510	1513	rBV4	4022	6186	0.25%	0.044%
18	12.443	1597	1601	1609	rBV4	3721	7223	0.29%	0.051%
19	13.571	1789	1793	1797	rVB2	2272	2870	0.11%	0.020%
20	14.094	1876	1882	1888	rBV	420852	624979	24.90%	4.407%
21	14.388	1924	1932	1942	rBV	457593	705615	28.12%	4.975%
22	14.693	1974	1984	1991	rBV	204780	320981	12.79%	2.263%
23	14.887	2007	2017	2029	rBV2	186912	323319	12.88%	2.280%
24	15.686	2143	2153	2160	rBV	565819	863094	34.39%	6.085%
25	15.803	2167	2173	2181	rBV	218319	326158	13.00%	2.300%
26	15.921	2188	2193	2197	rBV2	7848	11540	0.46%	0.081%
27	17.442	2447	2452	2459	rVV	245885	384667	15.33%	2.712%
28	17.542	2463	2469	2477	rVB2	585736	904895	36.06%	6.380%
29	18.100	2560	2564	2569	rBV4	5656	8060	0.32%	0.057%
30	18.153	2569	2573	2577	rVB3	3892	5097	0.20%	0.036%
31	18.329	2597	2603	2606	rBV4	15116	22426	0.89%	0.158%
32	18.758	2671	2676	2686	rBV	151288	226173	9.01%	1.595%
33	19.263	2758	2762	2766	rBV4	9087	12096	0.48%	0.085%
34	19.398	2780	2785	2788	rBV4	10312	12204	0.49%	0.086%
35	19.469	2793	2797	2800	rVV	8572	12361	0.49%	0.087%
36	19.510	2800	2804	2806	rVB4	2261	3302	0.13%	0.023%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049648.D
 Acq On : 5 Aug 2021 13:53
 Operator : CG/JU
 Sample : M3162-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG07

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-BG080421.M
 Title : SVOA CALIBRATION

37	19.604	2816	2820	2821	rBV4	3683	5113	0.20%	0.036%
38	19.663	2828	2830	2833	rBV3	3214	3083	0.12%	0.022%
39	19.833	2852	2859	2866	rVV	689165	1012173	40.33%	7.137%
40	19.886	2866	2868	2872	rVB3	2337	3211	0.13%	0.023%
41	20.056	2893	2897	2899	rBV4	2267	2997	0.12%	0.021%
42	20.356	2946	2948	2952	rVB4	3155	3407	0.14%	0.024%
43	20.579	2983	2986	2991	rVB5	4937	7490	0.30%	0.053%
44	20.620	2991	2993	2996	rBV3	2730	3304	0.13%	0.023%
45	20.726	3006	3011	3012	rBV3	4986	5693	0.23%	0.040%
46	20.802	3019	3024	3029	rBV	2080207	2509735	100.00%	17.696%
47	21.137	3078	3081	3085	rBV6	4049	5942	0.24%	0.042%
48	21.173	3085	3087	3090	rVV3	3082	3063	0.12%	0.022%
49	21.214	3090	3094	3098	rVB6	6364	8654	0.34%	0.061%
50	21.361	3112	3119	3124	rBV4	26860	54203	2.16%	0.382%
51	21.490	3140	3141	3146	rVB5	4918	5125	0.20%	0.036%
52	21.607	3160	3161	3165	rVB3	4798	3636	0.14%	0.026%
53	21.725	3174	3181	3187	rBV	239762	399003	15.90%	2.813%
54	21.907	3209	3212	3216	rBV5	5460	8153	0.32%	0.057%
55	22.441	3300	3303	3304	rBV3	4720	4573	0.18%	0.032%
56	22.500	3304	3313	3315	rBV7	19998	43710	1.74%	0.308%
57	23.152	3421	3424	3427	rBV4	3198	3565	0.14%	0.025%
58	23.193	3427	3431	3432	rVV3	2427	3441	0.14%	0.024%
59	23.234	3432	3438	3445	rVB7	8806	19420	0.77%	0.137%
60	23.375	3459	3462	3465	rBV3	2631	4188	0.17%	0.030%
61	23.428	3468	3471	3475	rVB4	3959	5557	0.22%	0.039%
62	23.699	3516	3517	3520	rBV3	3363	3718	0.15%	0.026%
63	23.734	3522	3523	3526	rVV3	4101	3488	0.14%	0.025%
64	23.757	3526	3527	3531	rVV4	3585	3408	0.14%	0.024%
65	23.910	3546	3553	3563	rVB7	33056	78874	3.14%	0.556%
66	24.057	3571	3578	3582	rVV5	12107	22225	0.89%	0.157%
67	24.779	3688	3701	3713	rBV3	342003	974571	38.83%	6.871%
68	24.915	3721	3724	3726	rBV3	3137	3417	0.14%	0.024%
69	25.003	3729	3739	3749	rVV2	142001	433724	17.28%	3.058%
70	25.138	3761	3762	3766	rVB4	4477	4628	0.18%	0.033%
71	25.408	3805	3808	3809	rBV3	3043	3242	0.13%	0.023%
72	25.437	3809	3813	3814	rBV3	8484	10445	0.42%	0.074%
73	25.531	3826	3829	3830	rBV3	3183	3754	0.15%	0.026%
74	25.755	3864	3867	3869	rVV3	4276	5650	0.23%	0.040%
75	25.831	3879	3880	3885	rVV4	4494	4013	0.16%	0.028%
76	26.166	3929	3937	3945	rBV7	13330	39410	1.57%	0.278%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
 Data File : BG049648.D
 Acq On : 5 Aug 2021 13:53
 Operator : CG/JU
 Sample : M3162-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGET7

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-BG080421.M
 Title : SVOA CALIBRATION

77	26.248	3949	3951	3954	rBV4	2344	2892	0.12%	0.020%
78	26.448	3984	3985	3990	rVV3	2398	2951	0.12%	0.021%
79	26.583	4005	4008	4012	rBV5	3257	5905	0.24%	0.042%
80	27.147	4101	4104	4107	rVB4	2569	3583	0.14%	0.025%
81	27.170	4107	4108	4111	rBV3	3140	3276	0.13%	0.023%
82	27.323	4130	4134	4137	rVB7	3319	4799	0.19%	0.034%
83	27.394	4145	4146	4153	rVB5	2768	2835	0.11%	0.020%
84	27.523	4166	4168	4169	rBV2	4534	3676	0.15%	0.026%
85	27.546	4169	4172	4174	rBV4	5124	5850	0.23%	0.041%
86	27.764	4205	4209	4212	rBV4	3396	5224	0.21%	0.037%
87	27.793	4212	4214	4218	rVB4	2244	2954	0.12%	0.021%
88	28.498	4332	4334	4337	rBV4	3071	3022	0.12%	0.021%
89	28.721	4369	4372	4375	rBV4	3247	5228	0.21%	0.037%
90	29.191	4450	4452	4454	rBV3	3863	4720	0.19%	0.033%
91	29.220	4454	4457	4458	rVV3	3712	4101	0.16%	0.029%
92	29.632	4524	4527	4528	rBV3	2664	3036	0.12%	0.021%
93	29.990	4586	4588	4591	rBV3	2628	3430	0.14%	0.024%
94	30.060	4596	4600	4603	rBV5	3136	4315	0.17%	0.030%
95	30.407	4655	4659	4661	rBV4	3570	4445	0.18%	0.031%
96	30.689	4705	4707	4711	rVB4	2897	3813	0.15%	0.027%
97	31.247	4797	4802	4803	rBV3	4430	7151	0.28%	0.050%
98	31.511	4844	4847	4850	rVB5	3634	4025	0.16%	0.028%
99	32.481	5009	5012	5014	rBV3	4386	4415	0.18%	0.031%
100	32.592	5027	5031	5034	rBV5	2041	2867	0.11%	0.020%

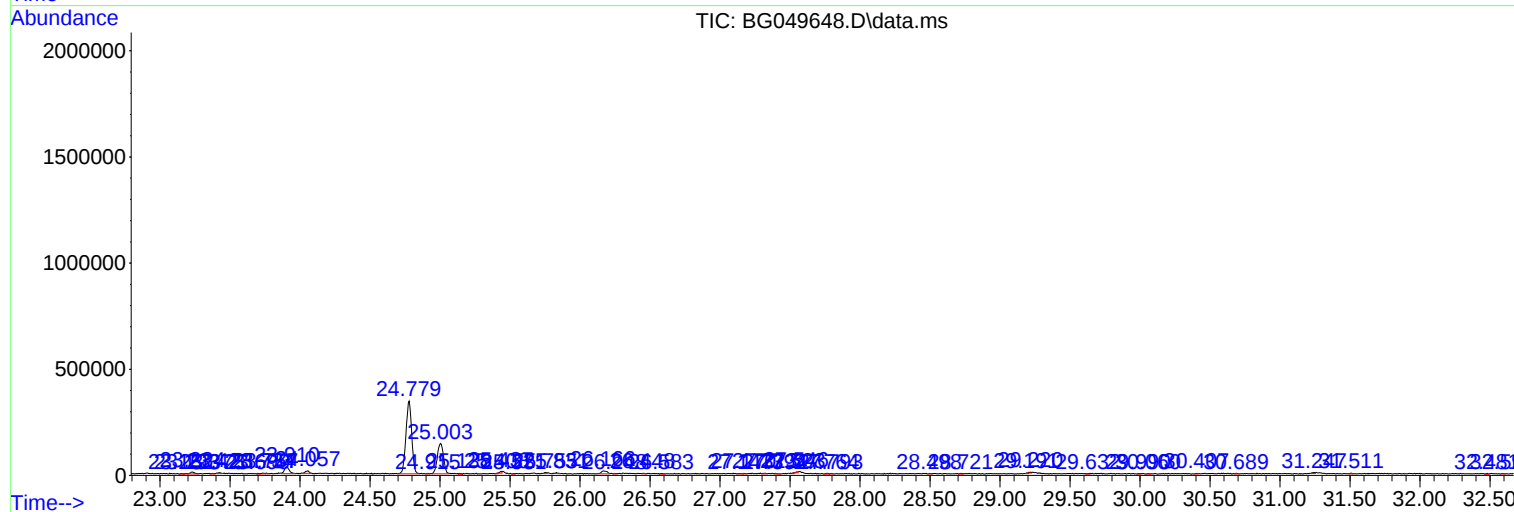
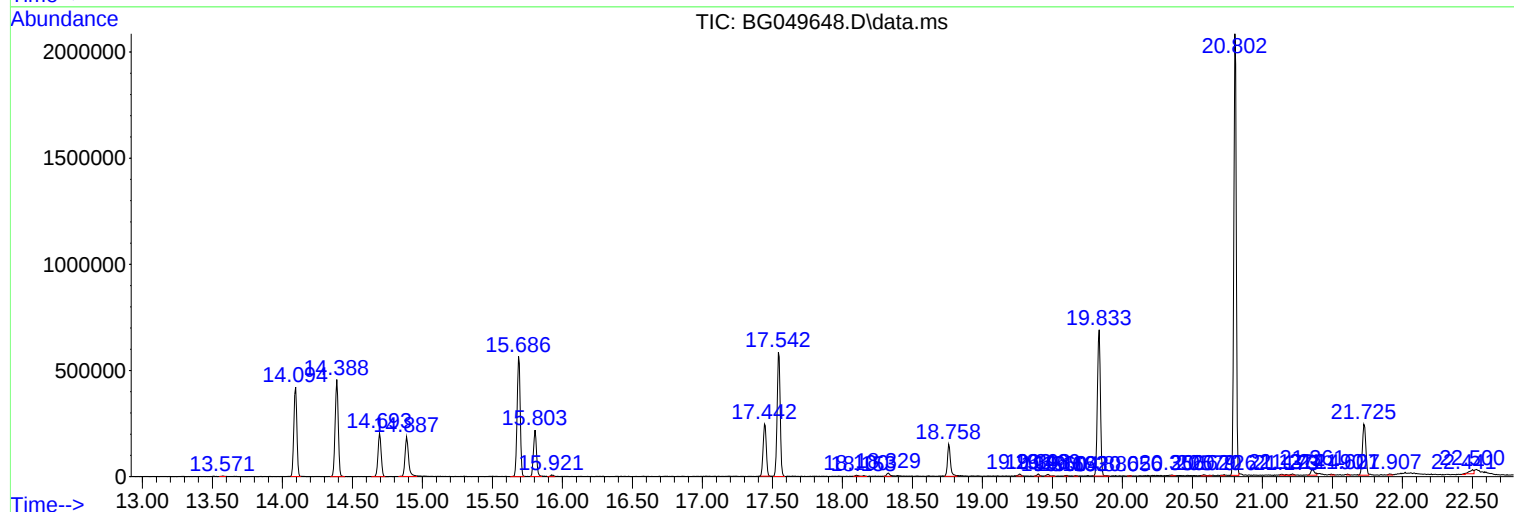
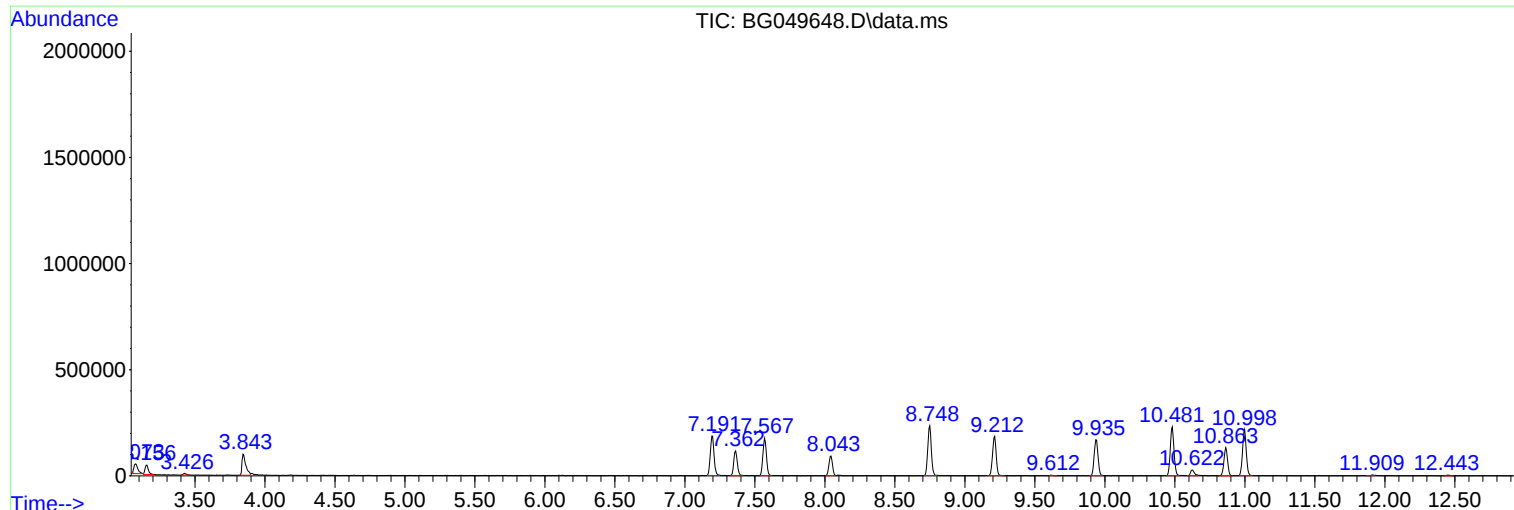
Sum of corrected areas: 14182853

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BG07

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

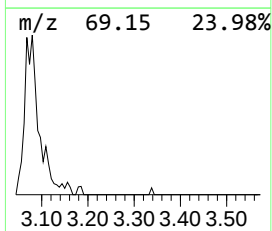
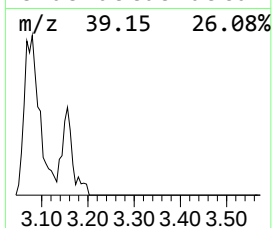
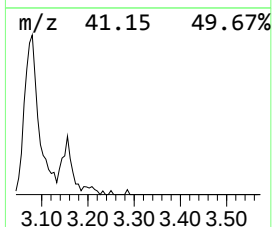
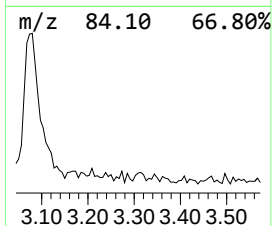
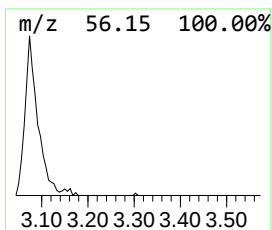
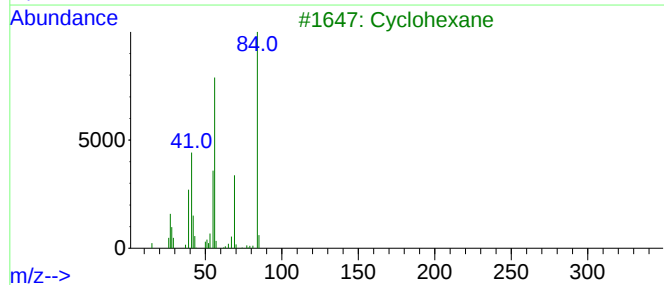
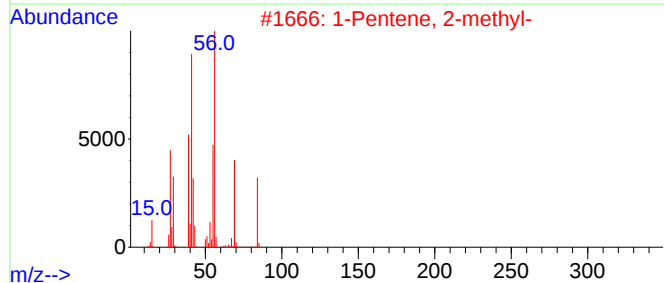
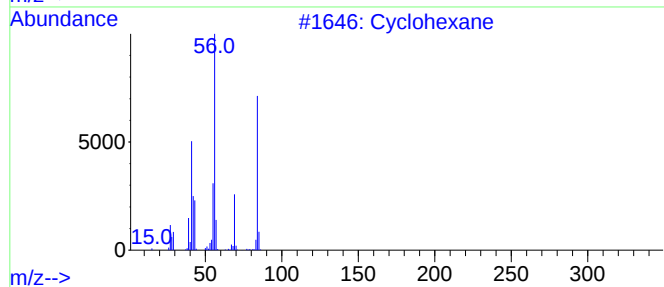
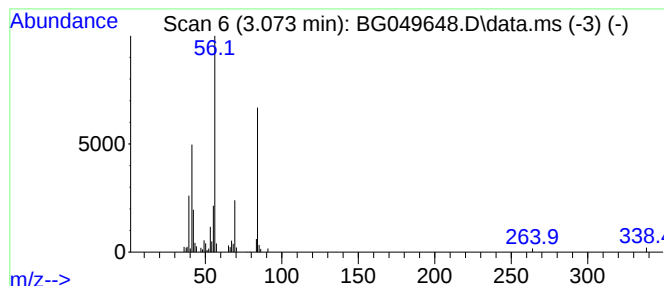
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.073 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.073	10.74 ng/ul	89127	1,4-Dichlorobenzene-d4	8.043

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	80
2			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	78
3			Cyclohexane	84	C6H12	000110-82-7	64
4			Cyclohexane	84	C6H12	000110-82-7	56
5			Iron pentacarbonyl	196	C5FeO5	013463-40-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BG07

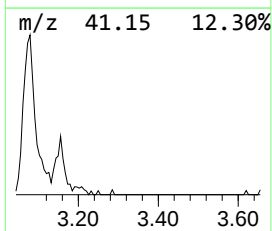
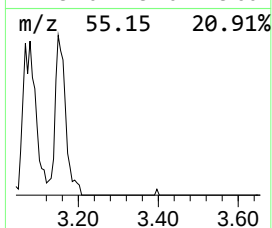
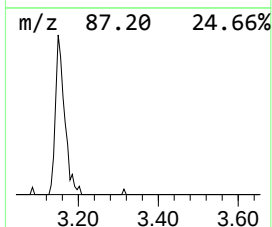
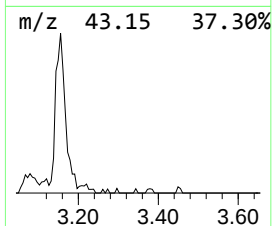
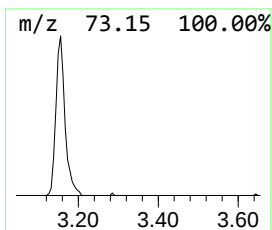
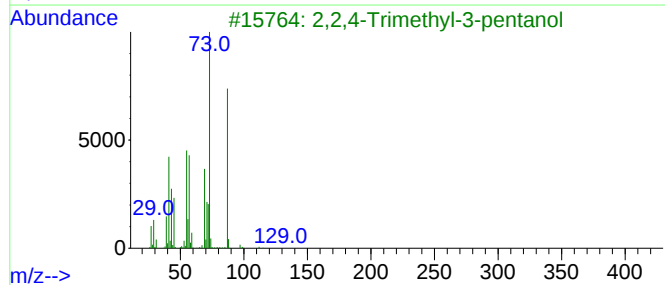
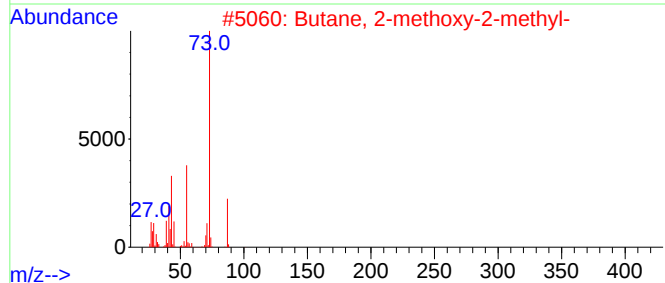
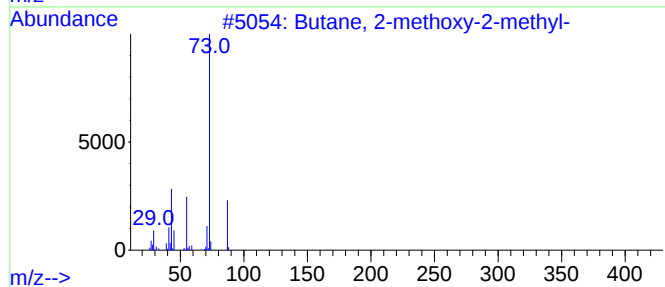
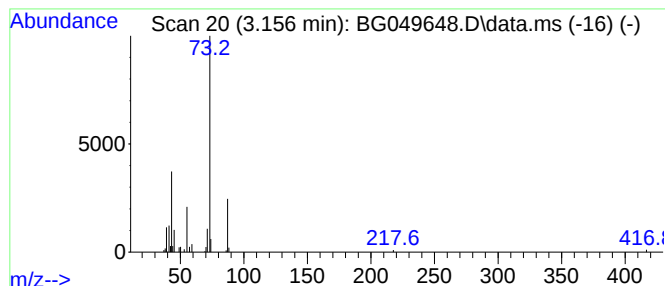
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.156	8.82 ng/ul	73212	1,4-Dichlorobenzene-d4	8.043

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
5			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

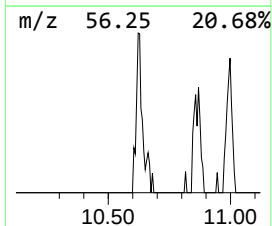
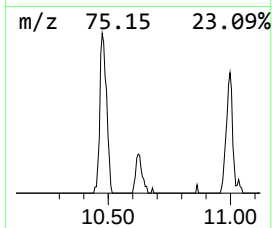
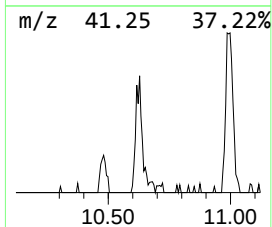
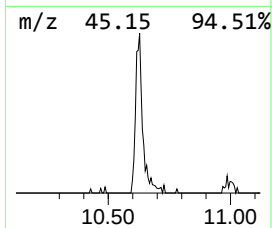
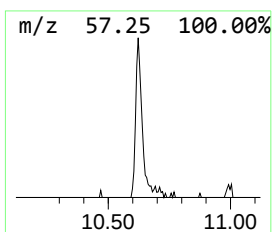
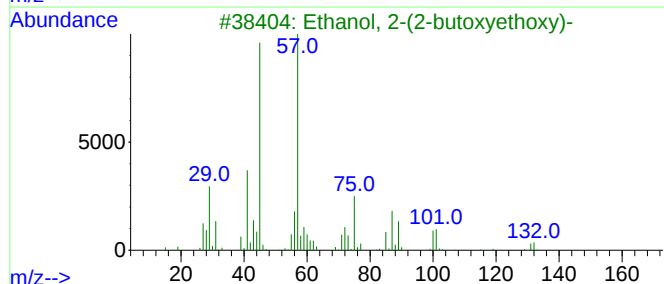
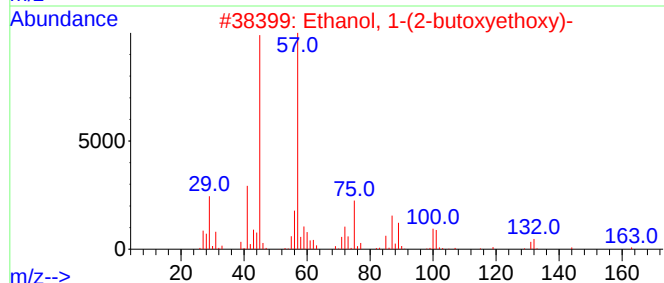
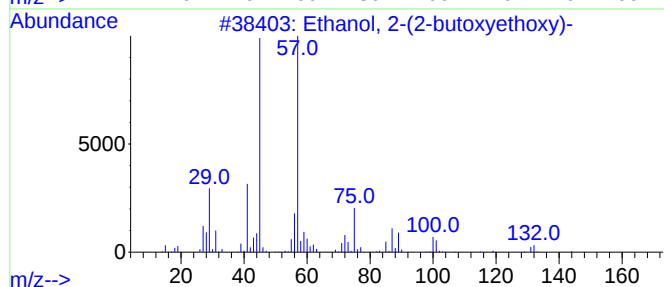
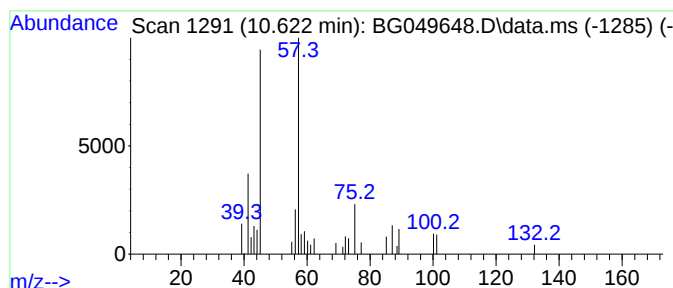
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.622	4.24 ng/ul	50609	Naphthalene-d8	10.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	86
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

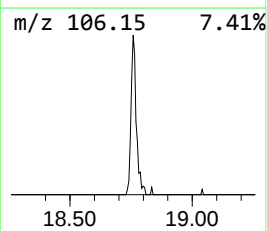
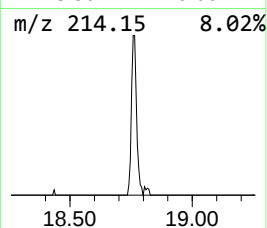
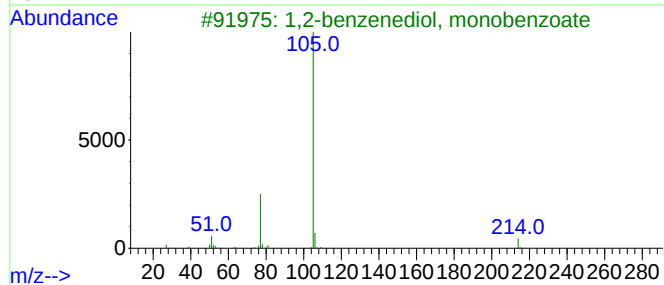
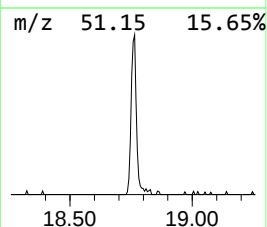
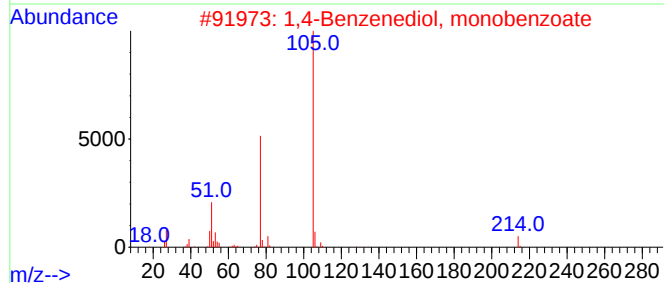
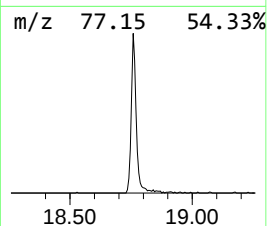
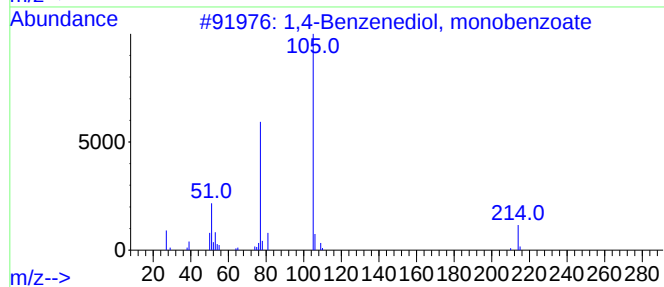
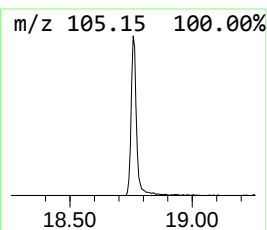
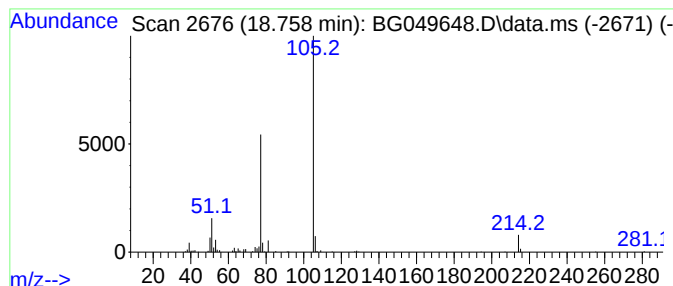
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,4-Benzenediol, monobenzoate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.758	11.76 ng/ul	226173	Phenanthrene-d10	17.442

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	91
2			1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	90
3			1,2-benzenediol, monobenzoate	214	C13H10O3	1000400-67-6	87
4			1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	83
5			N-(2,2-Dichlorovinyl)benzamide	215	C9H7Cl2NO	029431-43-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

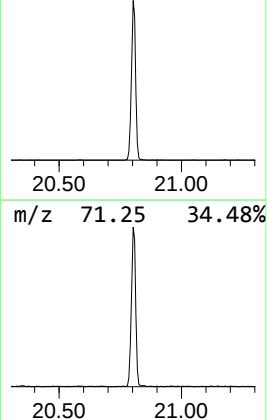
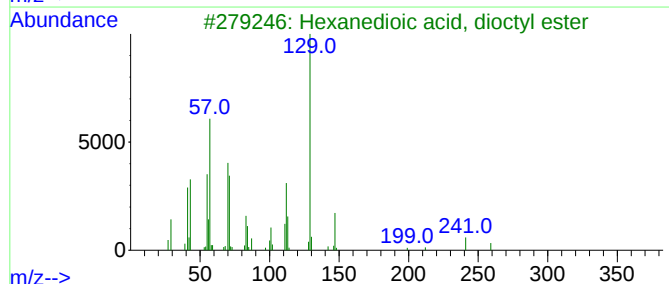
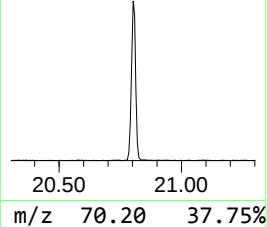
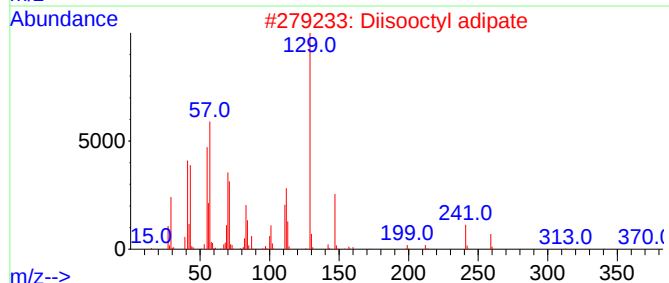
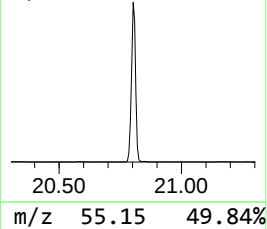
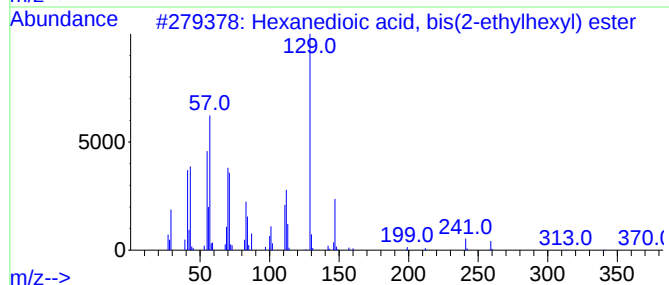
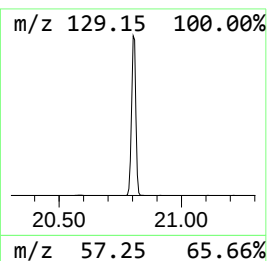
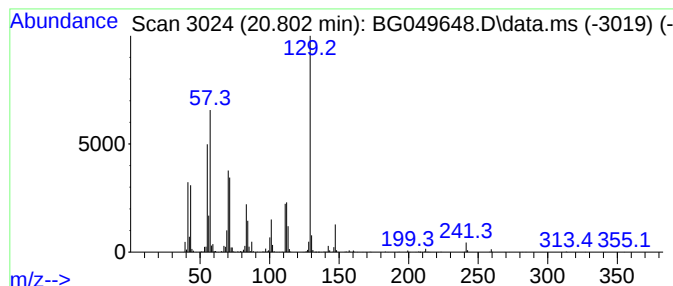
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexanedioic acid, bis(2-eth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.802	125.80 ng/ul	2509740	Chrysene-d12	21.725

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2			Diisooctyl adipate	370	C22H42O4	001330-86-5	91
3			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
4			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	81
5			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

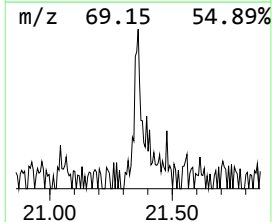
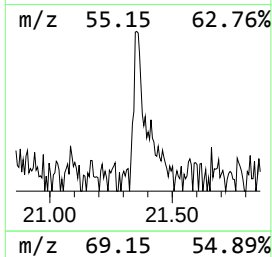
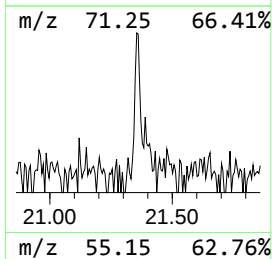
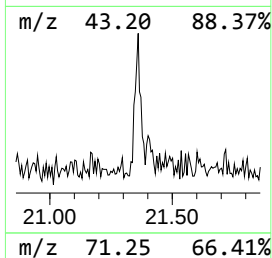
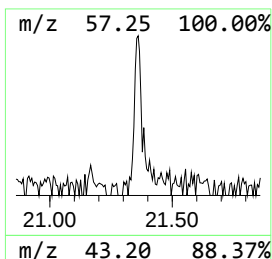
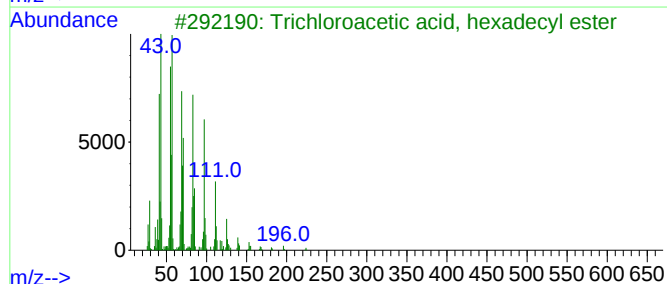
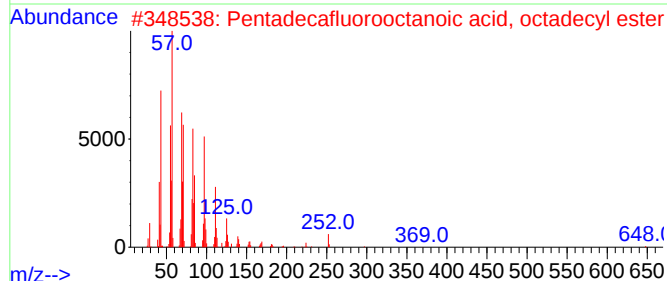
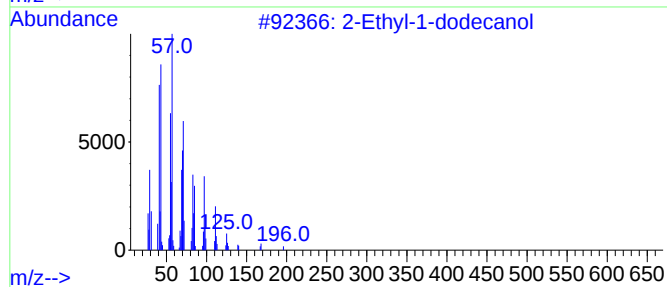
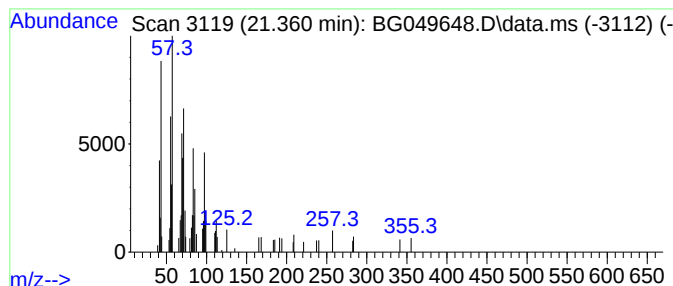
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Ethyl-1-dodecanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.360	2.72 ng/ul	54203	Chrysene-d12	21.725

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethyl-1-dodecanol	214	C14H30O	019780-33-7	72
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	62
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	53
4			Cyclododecanol, 1-ethenyl-	210	C14H26O	006244-49-1	49
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

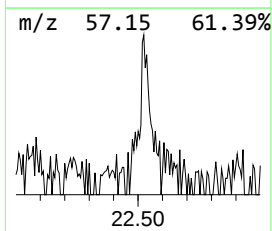
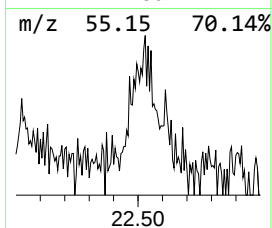
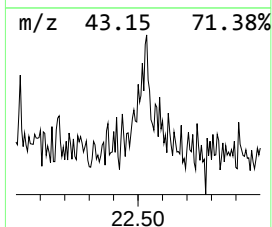
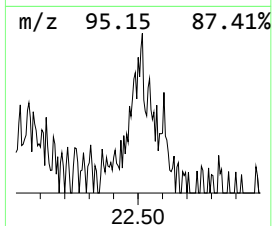
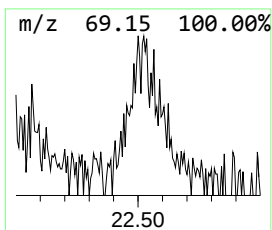
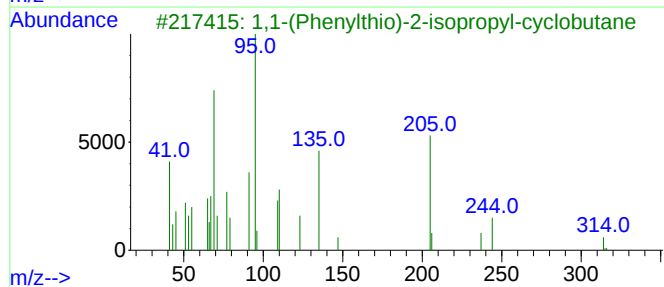
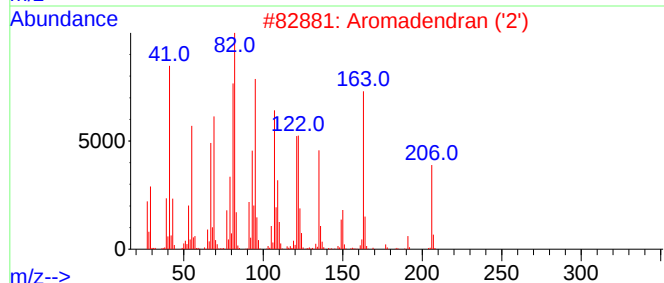
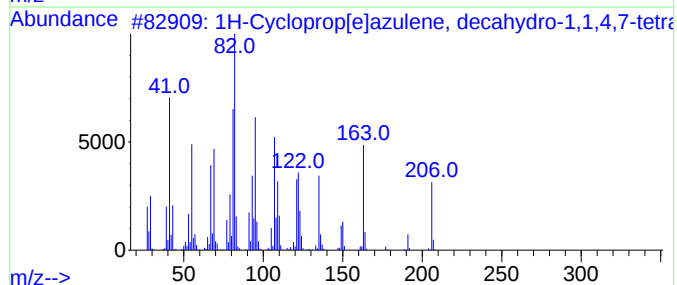
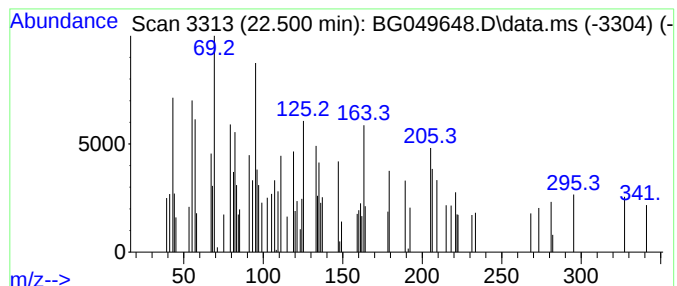
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.500	2.19 ng/ul	43710	Chrysene-d12	21.725

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cycloprop[e]azulene, decahydr...	206	C15H26	006790-78-9	27
2			Aromadendran ('2')	206	C15H26	110769-62-5	27
3			1,1-(Phenylthio)-2-isopropyl-cyc...	314	C19H22S2	122732-90-5	16
4			6,10-Dodecadien-3-ol, 3,7,11-tri...	224	C15H28O	003625-49-8	16
5			1-(2-Furyl)-4,4,4-trifluoro-1,3-...	206	C8H5F3O3	000326-90-9	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

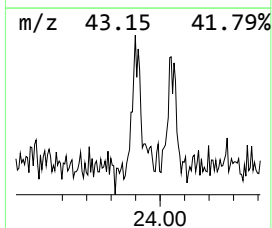
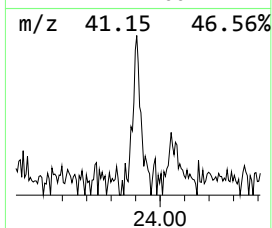
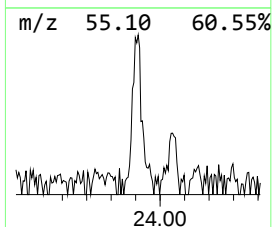
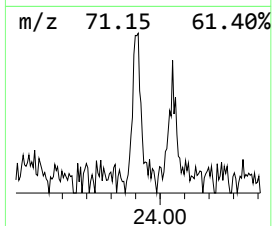
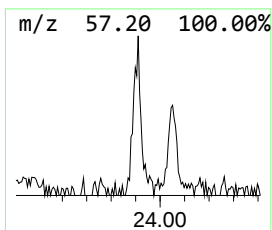
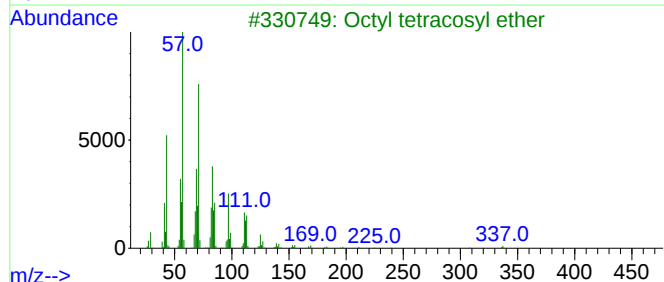
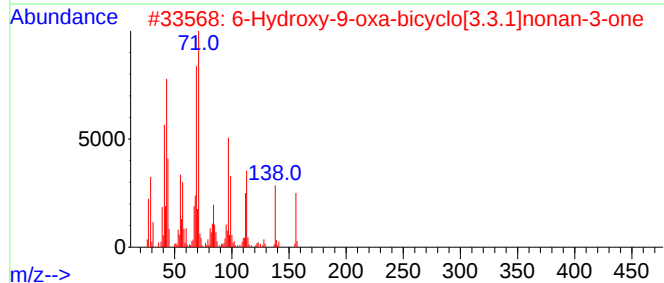
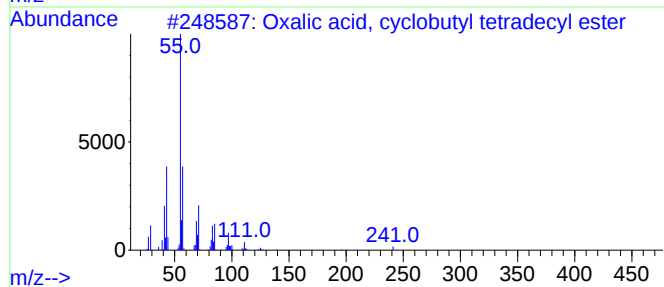
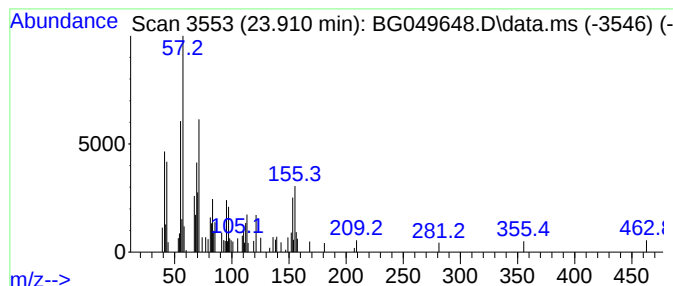
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.910	3.64 ng/ul	78874	Perylene-d12	25.003

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclobutyl tetradec...	340	C20H36O4	1000309-70-9	27
2			6-Hydroxy-9-oxa-bicyclo[3.3.1]no...	156	C8H12O3	1000194-46-8	25
3			Octyl tetracosyl ether	467	C32H66O	1000406-39-0	25
4			4-t-Pentylcyclohexene	152	C11H20	051874-62-5	25
5			Undecane	156	C11H24	001120-21-4	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049648.D
Acq On : 5 Aug 2021 13:53
Operator : CG/JU
Sample : M3162-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGET7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.073	10.7	ng/ul	89127	1	8.043	166022	20.0
Butane, 2-metho...	3.156	8.8	ng/ul	73212	1	8.043	166022	20.0
Ethanol, 2-(2-b...	10.622	4.2	ng/ul	50609	2	10.863	238791	20.0
1,4-Benzenediol...	18.758	11.8	ng/ul	226173	4	17.442	384667	20.0
Hexanedioic aci...	20.802	125.8	ng/ul	2509740	5	21.725	399003	20.0
2-Ethyl-1-dodec...	21.360	2.7	ng/ul	54203	5	21.725	399003	20.0
unknown-01	22.500	2.2	ng/ul	43710	5	21.725	399003	20.0
unknown-02	23.910	3.6	ng/ul	78874	6	25.003	433724	20.0