

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
 Data File : BG058207.D
 Acq On : 13 Jul 2023 20:52
 Operator : CG/JU
 Sample : 03414-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A46G0

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-BG070123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG058207.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	6	12	41	rVB	2404455	5021168	100.00%	29.127%
2	3.764	111	115	124	rBV	20515	36850	0.73%	0.214%
3	3.999	149	155	160	rBV	9058	13717	0.27%	0.080%
4	4.093	167	171	176	rVB3	6147	8793	0.18%	0.051%
5	4.339	205	213	221	rVB5	12464	25709	0.51%	0.149%
6	4.621	256	261	268	rVB2	16701	29324	0.58%	0.170%
7	5.003	320	326	332	rVB	6152	10489	0.21%	0.061%
8	5.368	383	388	390	rBV2	6062	8241	0.16%	0.048%
9	5.544	413	418	432	rVB2	10941	26469	0.53%	0.154%
10	5.802	452	462	474	rBV3	59169	151710	3.02%	0.880%
11	5.908	477	480	485	rBV4	4332	5836	0.12%	0.034%
12	6.184	520	527	534	rVB	26492	43656	0.87%	0.253%
13	6.525	578	585	598	rVB	157312	277438	5.53%	1.609%
14	7.066	669	677	691	rBV	96395	186844	3.72%	1.084%
15	7.230	698	705	714	rBV	78248	136135	2.71%	0.790%
16	7.436	733	740	750	rBV2	93267	190227	3.79%	1.103%
17	7.518	751	754	759	rVB4	3953	6545	0.13%	0.038%
18	7.618	766	771	775	rBV2	5048	8547	0.17%	0.050%
19	7.906	811	820	828	rBV	151884	271113	5.40%	1.573%
20	7.976	828	832	836	rVV	12228	20246	0.40%	0.117%
21	8.023	838	840	843	rVV3	4152	5815	0.12%	0.034%
22	8.076	843	849	856	rVV2	42162	77428	1.54%	0.449%
23	8.153	856	862	867	rVV2	57184	103881	2.07%	0.603%
24	8.217	867	873	883	rVV	430852	759788	15.13%	4.407%
25	8.611	932	940	946	rVV	93509	183838	3.66%	1.066%
26	8.664	946	949	955	rVV2	16480	28968	0.58%	0.168%
27	8.728	955	960	971	rVB3	24188	50667	1.01%	0.294%
28	8.899	983	989	1001	rVB2	20555	41757	0.83%	0.242%
29	9.063	1010	1017	1025	rVV	93312	173546	3.46%	1.007%
30	9.157	1029	1033	1036	rVV4	5002	8882	0.18%	0.052%
31	9.198	1036	1040	1049	rVB5	10458	22811	0.45%	0.132%
32	9.733	1127	1131	1134	rBV3	3882	6859	0.14%	0.040%
33	9.786	1134	1140	1151	rVB2	77864	143521	2.86%	0.833%
34	10.327	1224	1232	1250	rBV2	124291	273067	5.44%	1.584%
35	10.708	1285	1297	1310	rVB	241704	456353	9.09%	2.647%
36	10.844	1314	1320	1330	rBV	81882	161297	3.21%	0.936%

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37	11.243	1383	1388	1395	rVB4	3164	6055	0.12%	0.035%
38	11.519	1426	1435	1442	rVB6	3903	10045	0.20%	0.058%
39	12.659	1624	1629	1633	rBV2	6394	10336	0.21%	0.060%
40	12.759	1641	1646	1648	rBV3	3102	5127	0.10%	0.030%

41	12.782	1648	1650	1659	rVB3	3565	6385	0.13%	0.037%
42	13.358	1744	1748	1752	rVB2	5149	7053	0.14%	0.041%
43	13.429	1752	1760	1765	rVB2	7845	13003	0.26%	0.075%
44	13.940	1841	1847	1856	rBV	246088	363663	7.24%	2.110%
45	14.181	1878	1888	1891	rBV3	16874	31924	0.64%	0.185%

46	14.234	1891	1897	1905	rVB	289751	456302	9.09%	2.647%
47	14.539	1942	1949	1959	rBV	430919	666544	13.27%	3.867%
48	14.745	1978	1984	1996	rBV2	57549	133786	2.66%	0.776%
49	14.892	2005	2009	2012	rVB4	5519	7132	0.14%	0.041%
50	15.532	2112	2118	2128	rBV	381129	578556	11.52%	3.356%

51	15.650	2131	2138	2151	rVB	83216	132801	2.64%	0.770%
52	15.896	2174	2180	2186	rBV2	10802	16999	0.34%	0.099%
53	16.073	2206	2210	2218	rBV5	6369	15182	0.30%	0.088%
54	16.972	2359	2363	2370	rVB	53050	71508	1.42%	0.415%
55	17.283	2410	2416	2426	rBV	566379	866917	17.27%	5.029%

56	17.383	2427	2433	2445	rVV	393057	599148	11.93%	3.476%
57	17.941	2525	2528	2535	rVB3	10810	13842	0.28%	0.080%
58	18.170	2563	2567	2575	rBV4	12221	24658	0.49%	0.143%
59	18.235	2576	2578	2583	rVV2	3631	6902	0.14%	0.040%
60	18.717	2656	2660	2662	rVV2	3462	5537	0.11%	0.032%

61	18.746	2662	2665	2671	rVB3	5292	7985	0.16%	0.046%
62	18.899	2688	2691	2696	rVV5	3551	6880	0.14%	0.040%
63	18.975	2698	2704	2708	rVB4	8437	11349	0.23%	0.066%
64	19.022	2708	2712	2716	rBV4	4205	6541	0.13%	0.038%
65	19.110	2722	2727	2731	rBV6	10798	14454	0.29%	0.084%

66	19.245	2745	2750	2757	rVB5	9398	18893	0.38%	0.110%
67	19.451	2779	2785	2787	rBV7	4584	7020	0.14%	0.041%
68	19.586	2806	2808	2813	rVB5	4730	5858	0.12%	0.034%
69	19.674	2816	2823	2835	rBV	541204	782889	15.59%	4.541%
70	19.898	2858	2861	2864	rVB2	7341	7950	0.16%	0.046%

71	20.139	2899	2902	2906	rBV5	5594	7961	0.16%	0.046%
72	20.191	2909	2911	2917	rVB6	5673	8802	0.18%	0.051%
73	20.691	2993	2996	3000	rVB5	9115	11866	0.24%	0.069%
74	20.908	3031	3033	3037	rVB4	4966	6284	0.13%	0.036%
75	20.990	3042	3047	3050	rBV2	25417	35100	0.70%	0.204%

76	21.184	3078	3080	3084	rBV4	6692	7902	0.16%	0.046%
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77	21.419	3117	3120	3124	rVB	24822	31002	0.62%	0.180%
78	21.537	3133	3140	3149	rBV2	560498	957721	19.07%	5.556%
79	21.719	3168	3171	3177	rVB	15567	19018	0.38%	0.110%
80	22.307	3267	3271	3280	rVB	20942	33703	0.67%	0.196%
81	22.959	3376	3382	3395	rBV5	90125	210927	4.20%	1.224%
82	23.764	3512	3519	3523	rBV3	23387	49879	0.99%	0.289%
83	24.457	3627	3637	3648	rBV2	268545	741364	14.76%	4.301%
84	24.680	3664	3675	3688	rBV	382164	1090659	21.72%	6.327%
85	25.773	3855	3861	3870	rVB4	24631	64101	1.28%	0.372%
86	27.083	4078	4084	4095	rVB5	21731	65626	1.31%	0.381%

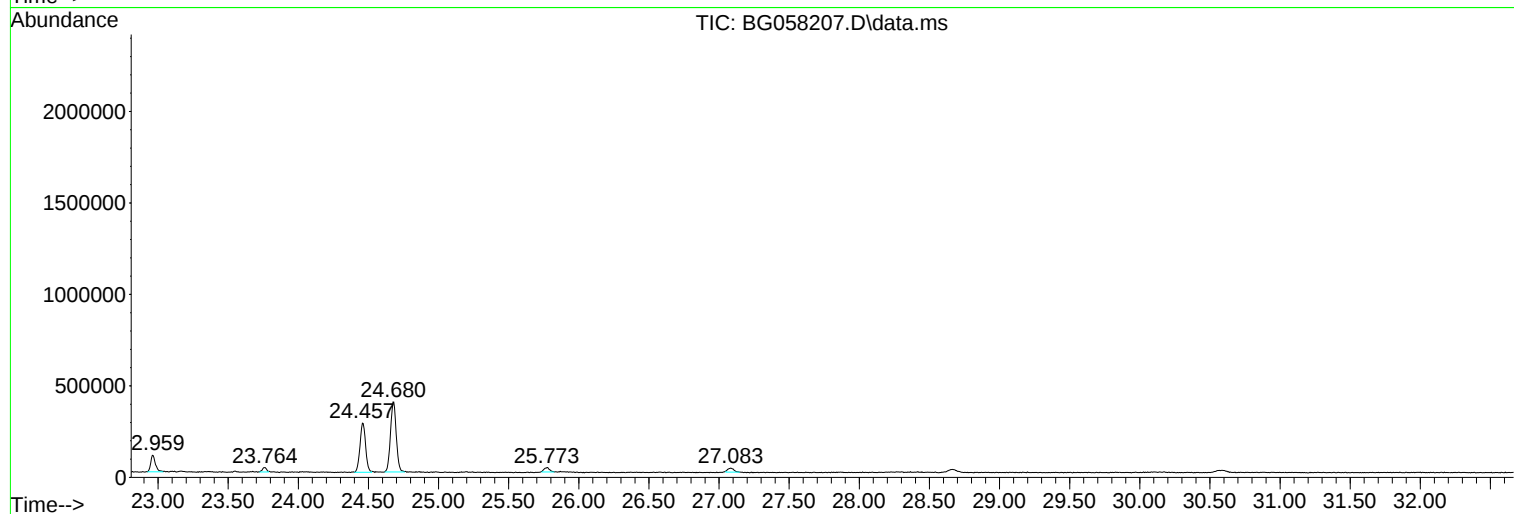
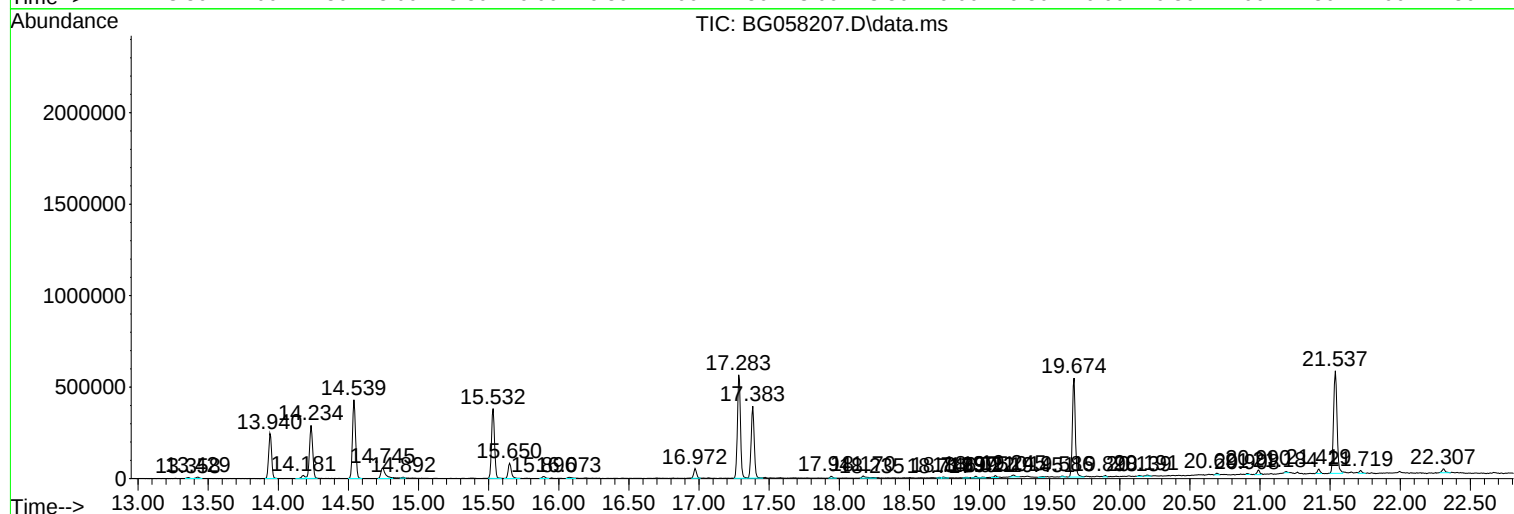
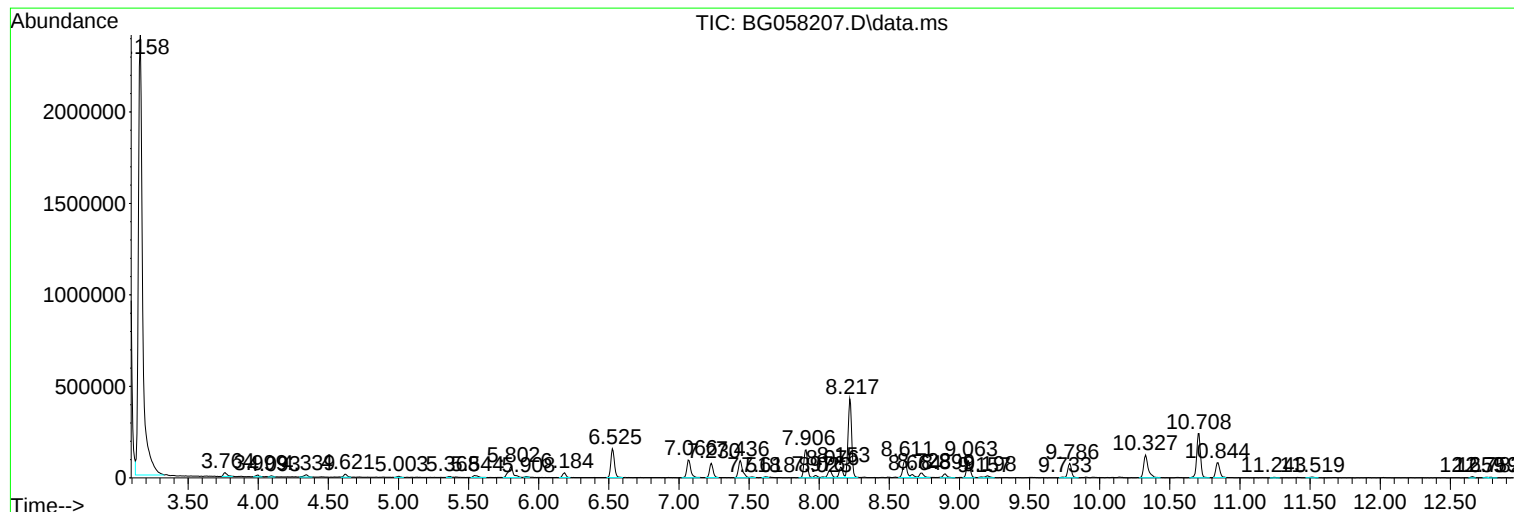
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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



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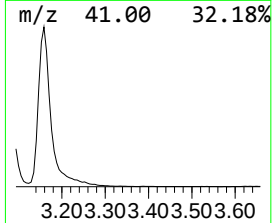
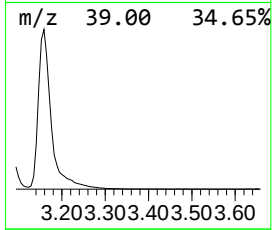
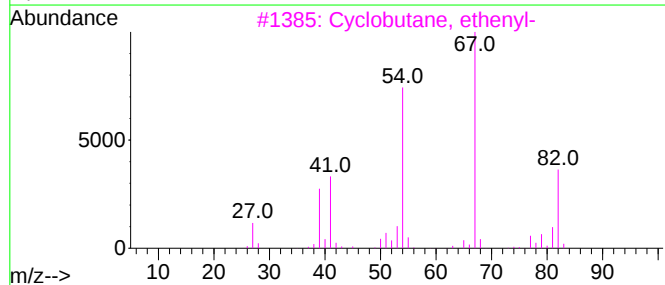
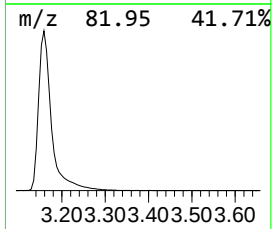
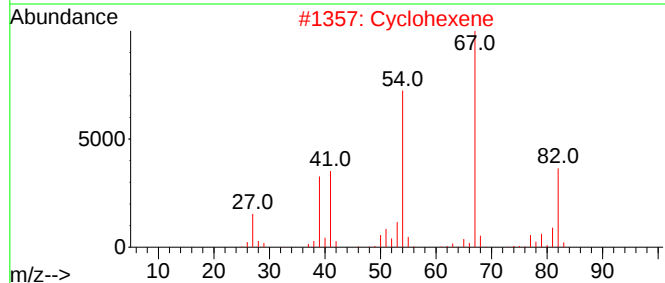
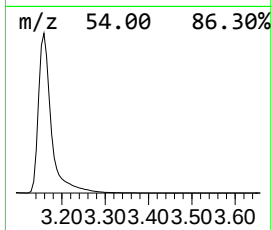
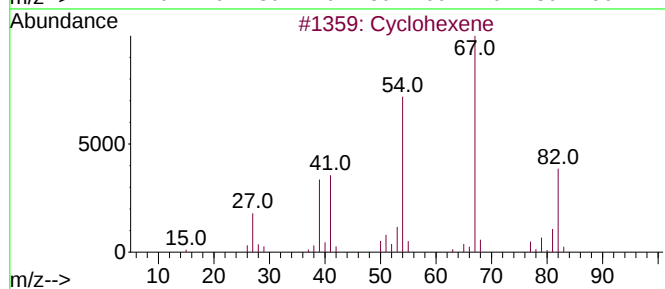
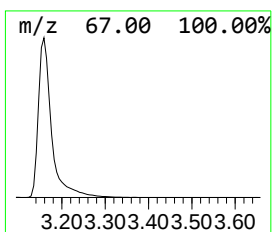
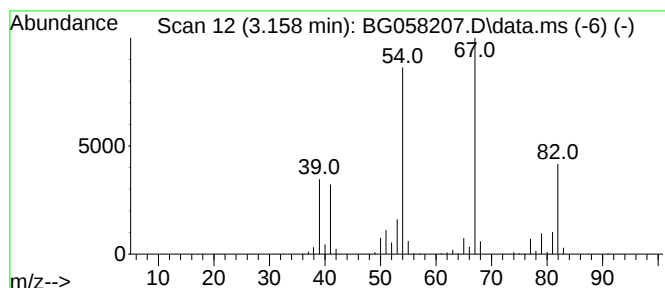
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclohexene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.158	370.41 ng/ul	5021170	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene	82	C6H10	000110-83-8	95
2			Cyclohexene	82	C6H10	000110-83-8	91
3			Cyclobutane, ethenyl-	82	C6H10	002597-49-1	91
4			Cyclohexene	82	C6H10	000110-83-8	91
5			Ethylidenecyclobutane	82	C6H10	001528-21-8	91



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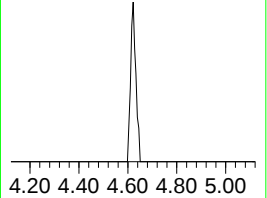
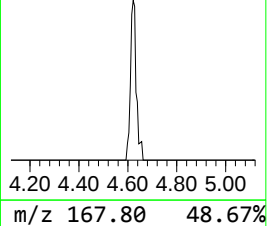
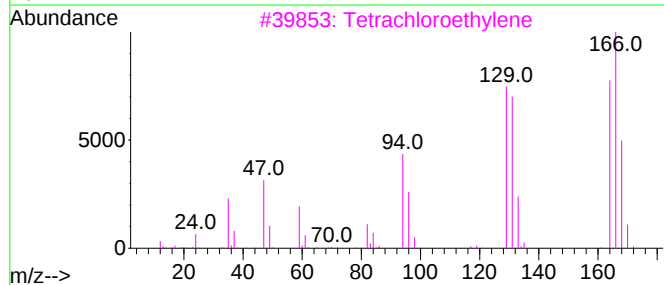
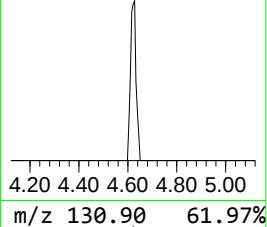
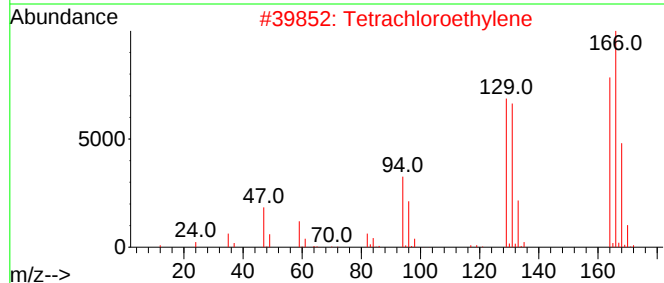
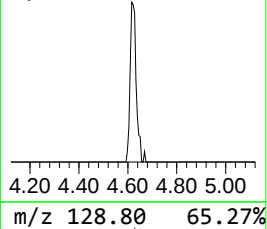
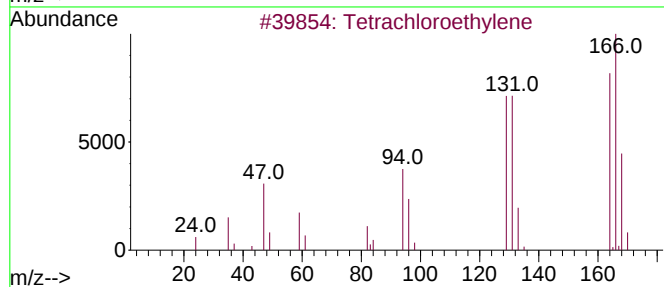
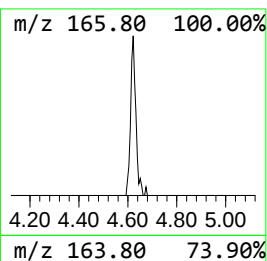
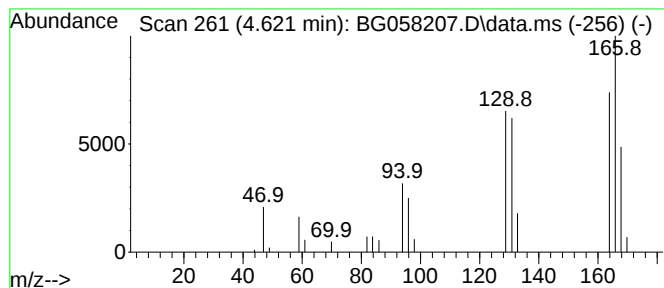
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Tetrachloroethylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.621	2.16 ng/ul	29324	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	94
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	72
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	59
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	43



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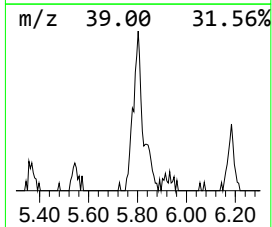
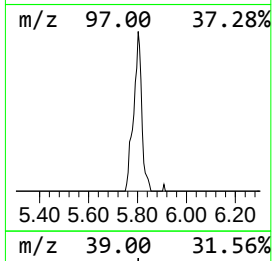
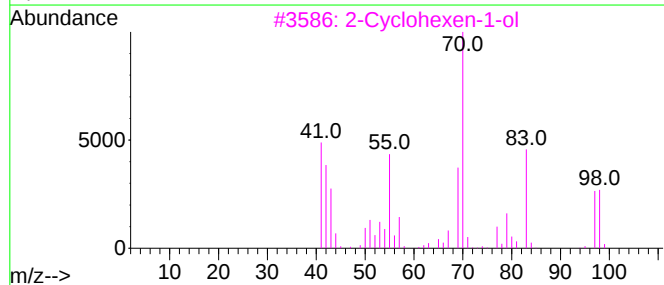
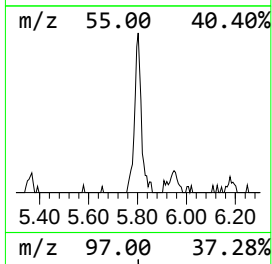
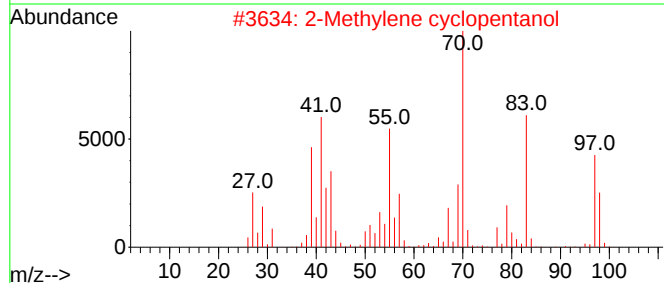
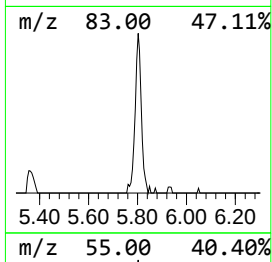
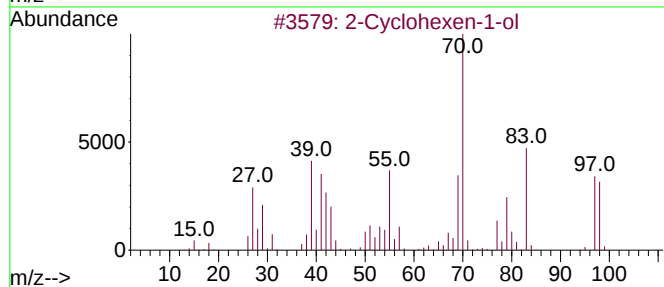
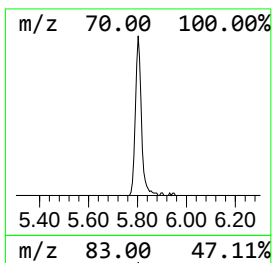
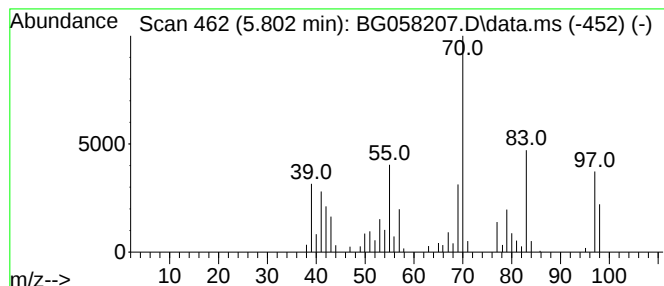
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Cyclohexen-1-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.802	11.19 ng/ul	151710	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	90
2	2-Methylene cyclopentanol			98	C6H10O	020461-31-8	90
3	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	87
4	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	81
5	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	59



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A46G0

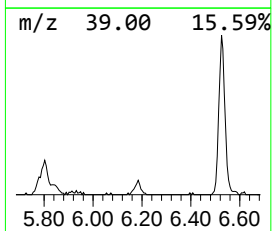
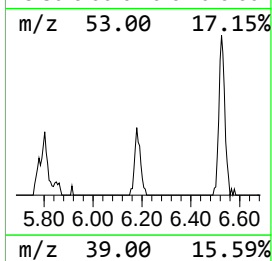
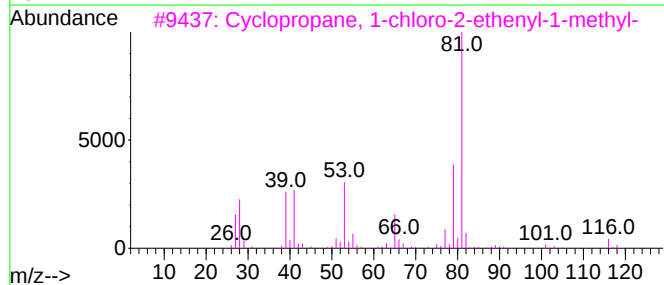
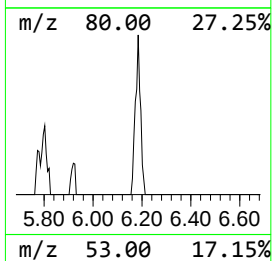
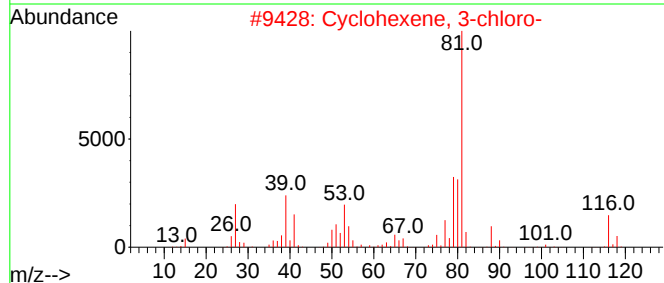
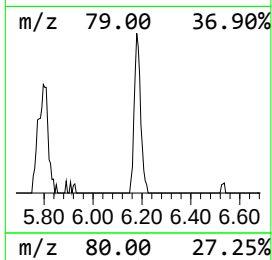
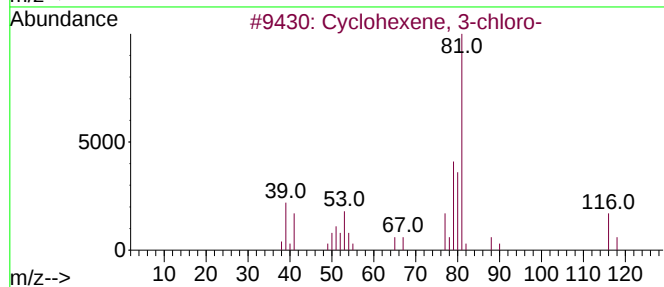
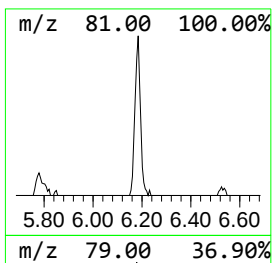
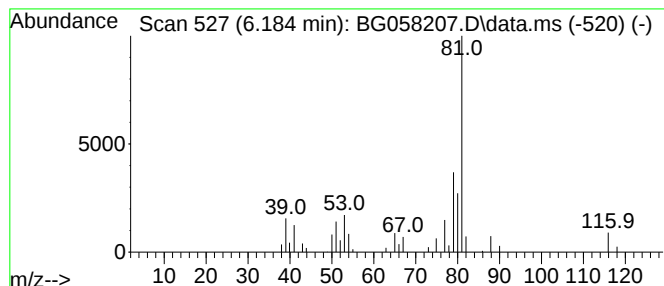
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclohexene, 3-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.184	3.22 ng/ul	43656	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 3-chloro-	116	C6H9Cl	002441-97-6	83
2			Cyclohexene, 3-chloro-	116	C6H9Cl	002441-97-6	74
3			Cyclopropane, 1-chloro-2-ethenyl...	116	C6H9Cl	062337-93-3	64
4			Cyclohexene, 1-chloro-	116	C6H9Cl	000930-66-5	59
5			Cyclohexene, 3-bromo-	160	C6H9Br	001521-51-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058207.D
Acq On : 13 Jul 2023 20:52
Operator : CG/JU
Sample : 03414-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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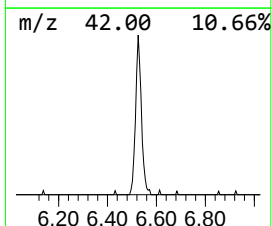
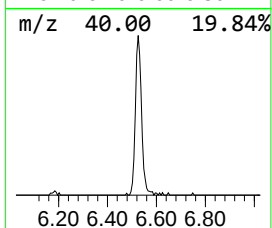
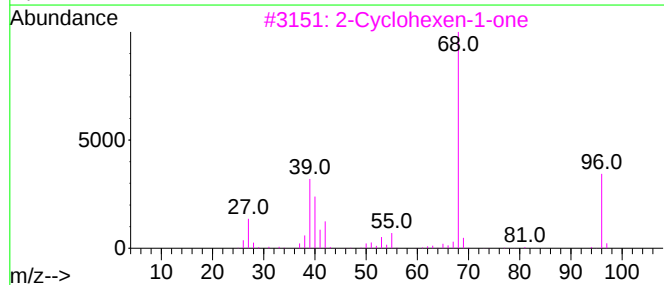
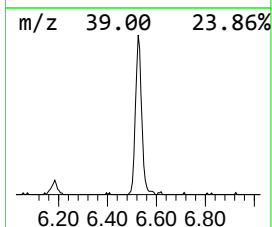
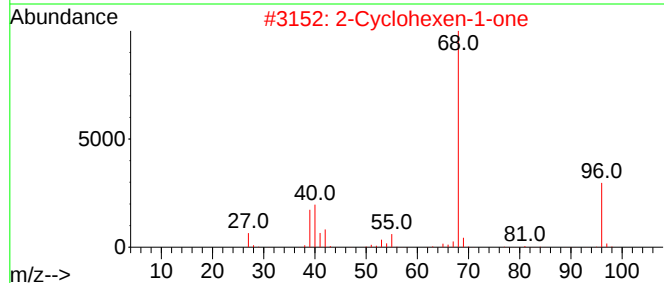
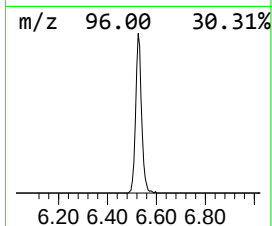
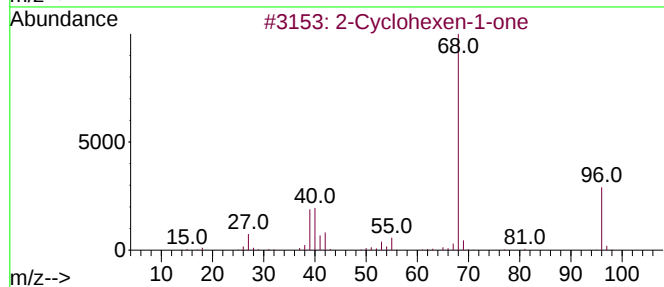
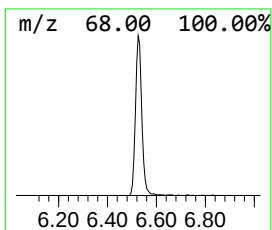
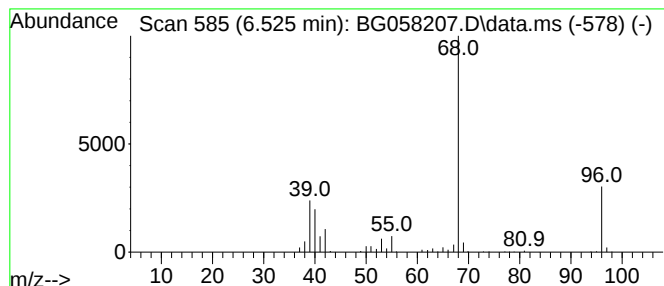
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Cyclohexen-1-one Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.525	20.47 ng/ul	277438	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	91
2			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	91
3			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	91
4			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	90
5			2H-Pyran-2-one, 5,6-dihydro-	98	C5H6O2	003393-45-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058207.D
Acq On : 13 Jul 2023 20:52
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Sample : 03414-06
Misc :
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Instrument :
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ClientSampleId :
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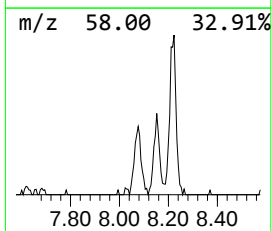
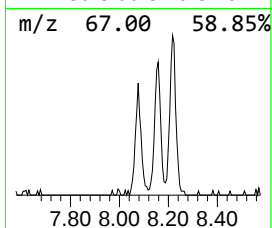
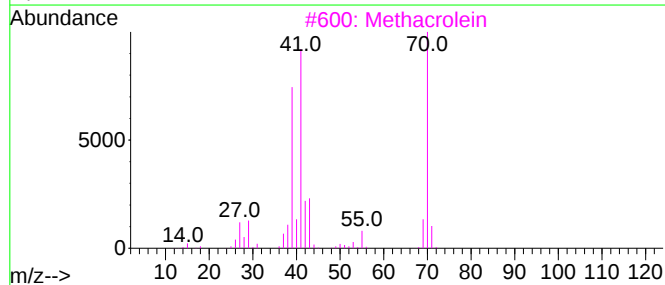
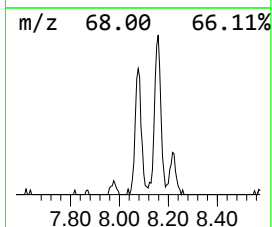
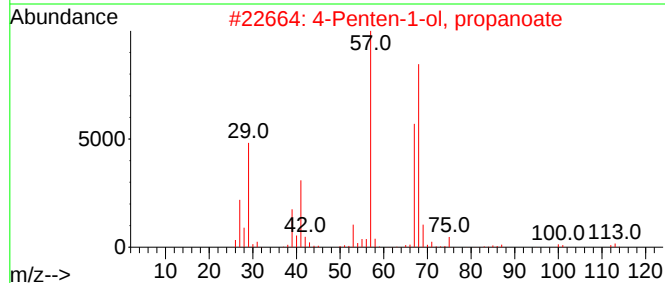
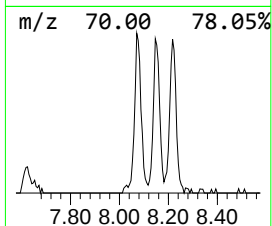
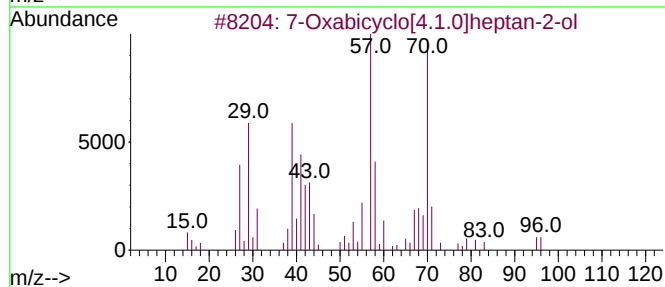
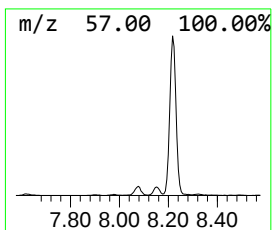
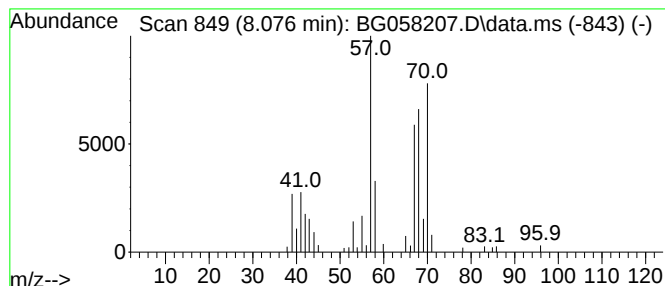
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.076	5.71 ng/ul	77428	1,4-Dichlorobenzene-d4	7.900

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		7-Oxabicyclo[4.1.0]heptan-2-ol	114	C6H10O2	001192-78-5	40
2		4-Penten-1-ol, propanoate	142	C8H14O2	030563-30-5	32
3		Methacrolein	70	C4H6O	000078-85-3	30
4		Ethyl pent-4-enyl carbonate	158	C8H14O3	1000373-78-4	17
5		cis,cis-2,7-Nonadiene	124	C9H16	036901-84-5	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058207.D
Acq On : 13 Jul 2023 20:52
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Sample : 03414-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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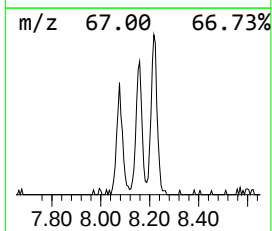
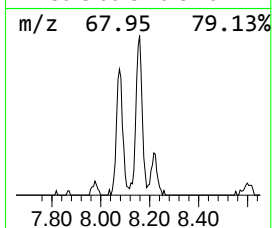
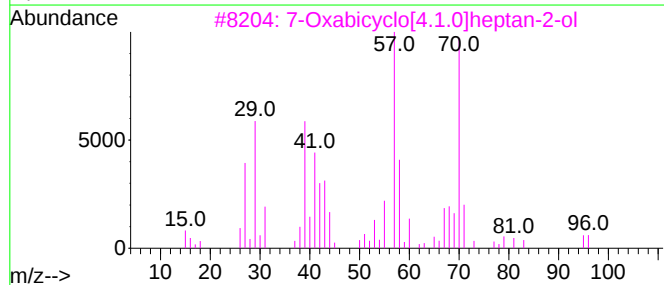
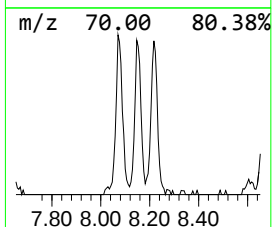
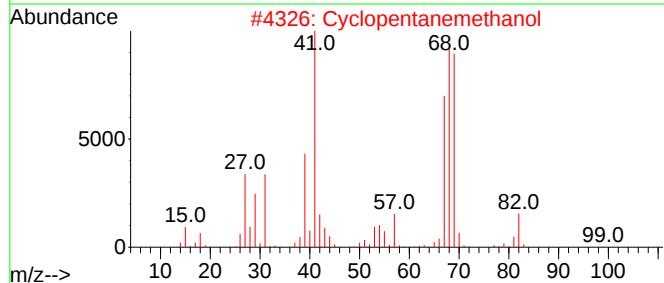
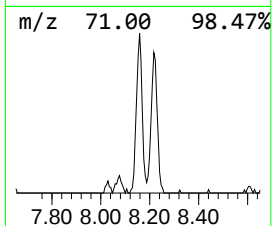
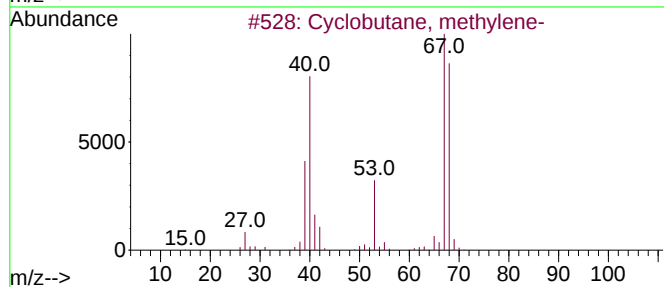
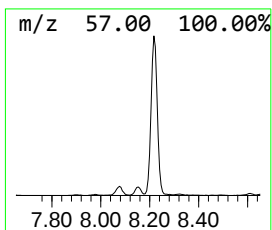
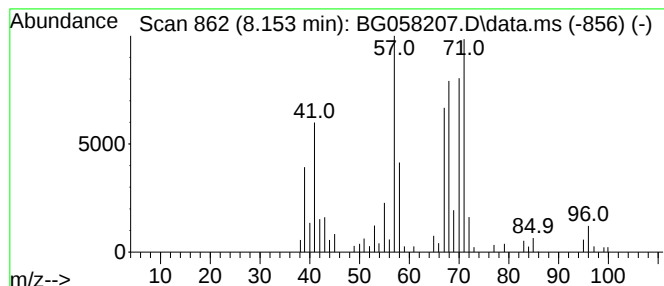
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.153	7.66 ng/ul	103881	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobutane, methylene-	68	C5H8	001120-56-5	27
2			Cyclopentanemethanol	100	C6H12O	003637-61-4	27
3			7-Oxabicyclo[4.1.0]heptan-2-ol	114	C6H10O2	001192-78-5	25
4			Oxirane, butyl-	100	C6H12O	001436-34-6	25
5			Cyclopentanol, 3-methyl-	100	C6H12O	018729-48-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058207.D
Acq On : 13 Jul 2023 20:52
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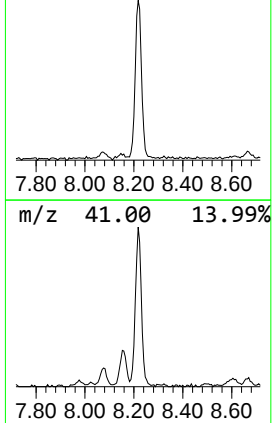
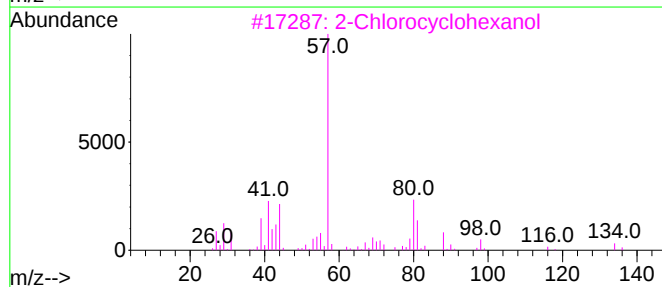
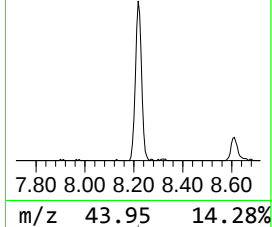
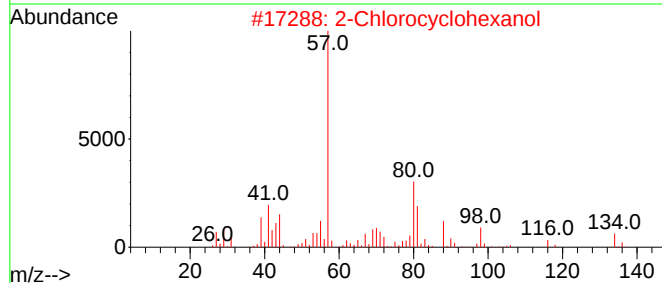
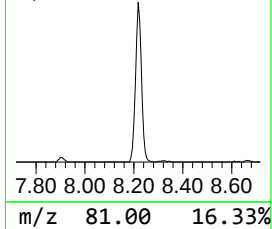
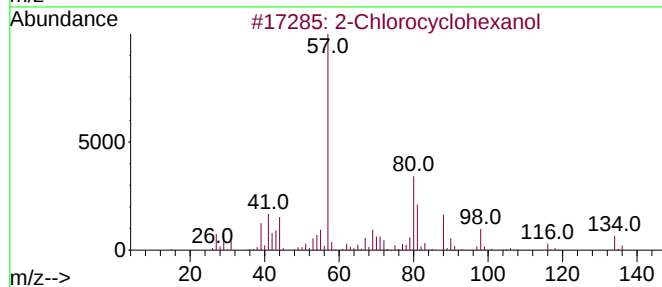
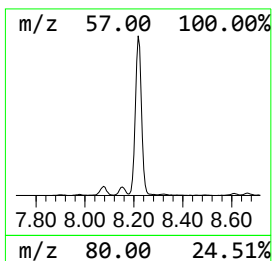
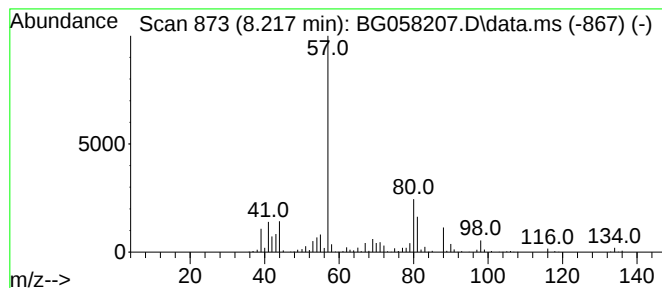
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Chlorocyclohexanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.217	56.05 ng/ul	759788	1,4-Dichlorobenzene-d4	7.900

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	95
2		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	94
3		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	91
4		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	83
5		2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058207.D
Acq On : 13 Jul 2023 20:52
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Sample : 03414-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

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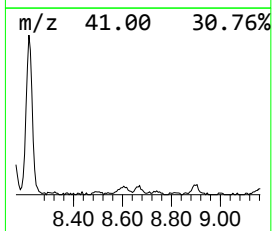
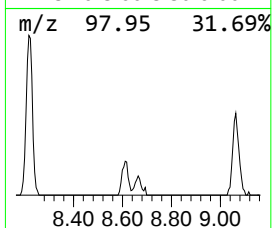
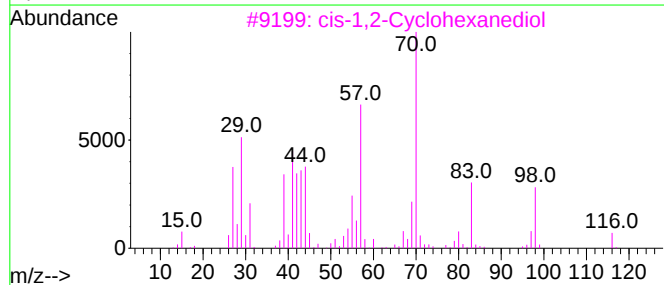
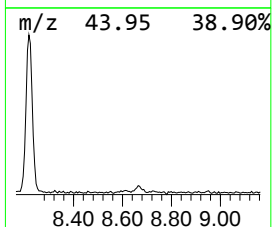
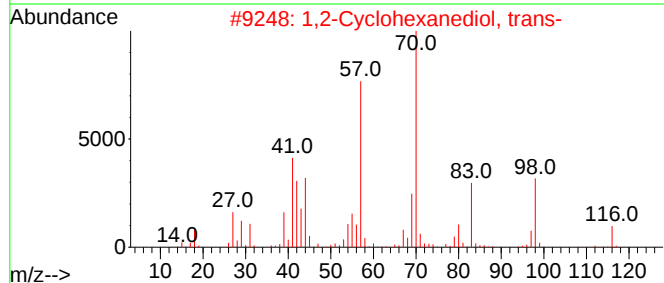
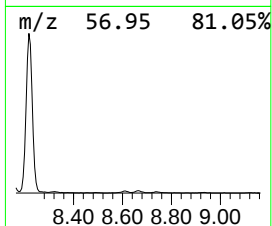
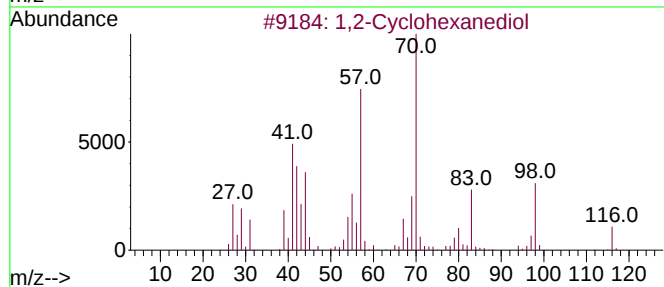
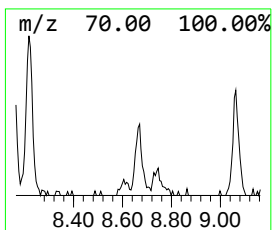
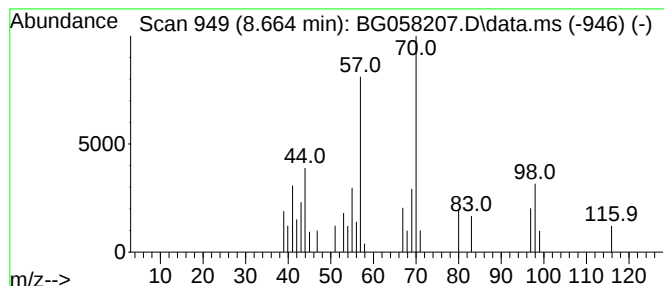
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,2-Cyclohexanediol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.664	2.14 ng/ul	28968	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Cyclohexanediol			116	C6H12O2	000931-17-9	72
2	1,2-Cyclohexanediol, trans-			116	C6H12O2	001460-57-7	64
3	cis-1,2-Cyclohexanediol			116	C6H12O2	001792-81-0	58
4	1,2-Cyclohexanediol			116	C6H12O2	000931-17-9	58
5	1,2-Cyclohexanediol, trans-			116	C6H12O2	001460-57-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
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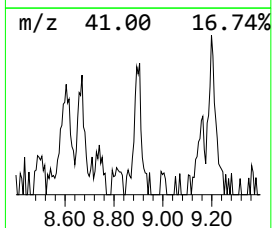
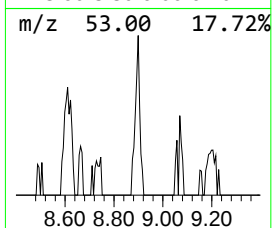
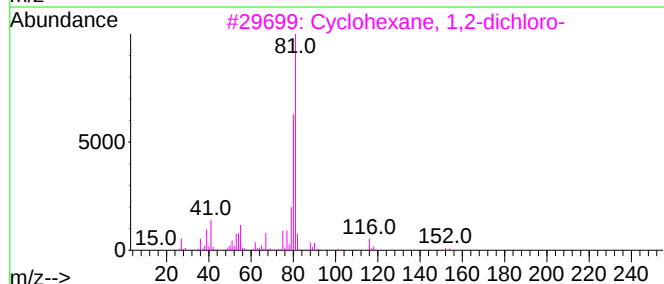
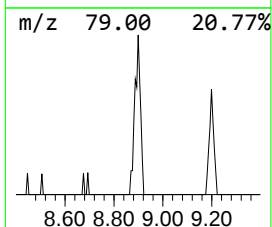
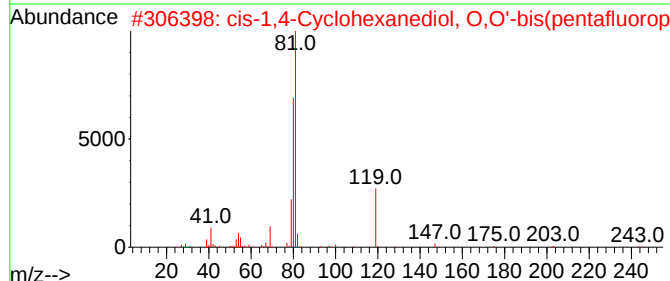
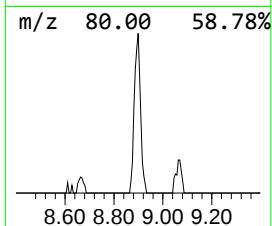
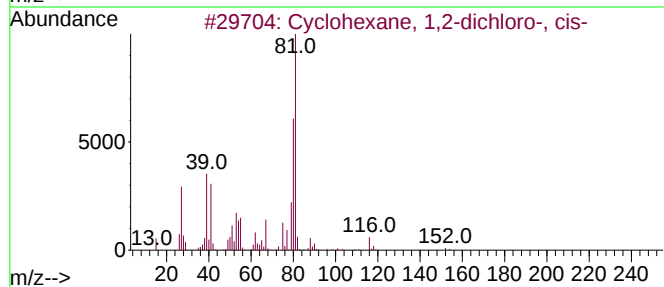
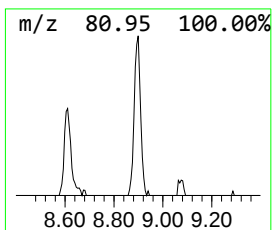
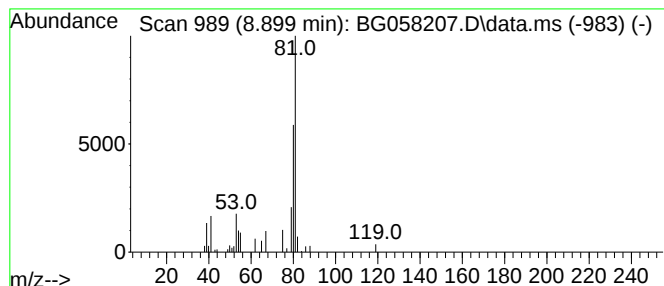
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Cyclohexane, 1,2-dichloro-,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.899	3.08 ng/ul	41757	1,4-Dichlorobenzene-d4	7.900

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dichloro-, cis-	152	C6H10Cl2	010498-35-8	74
2			cis-1,4-Cyclohexanediol, O,O'-bi...	408	C12H10F10O4	1000385-95-0	72
3			Cyclohexane, 1,2-dichloro-	152	C6H10Cl2	001121-21-7	64
4			Bi-2-cyclohexen-1-yl	162	C12H18	001541-20-4	64
5			Cyclohexane, 1,2-dichloro-, trans-	152	C6H10Cl2	000822-86-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058207.D
Acq On : 13 Jul 2023 20:52
Operator : CG/JU
Sample : 03414-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A46G0

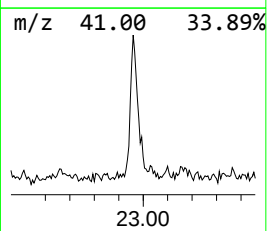
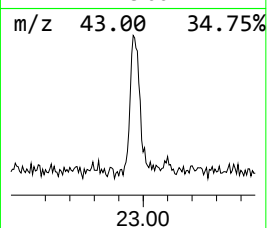
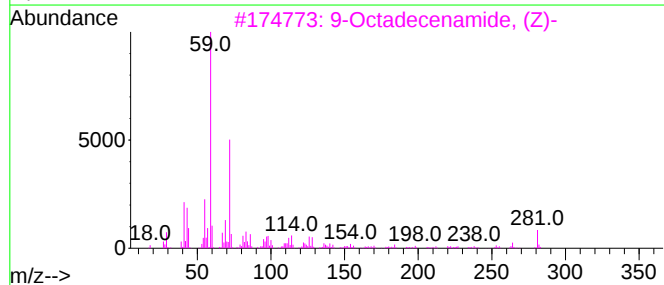
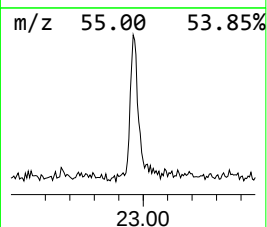
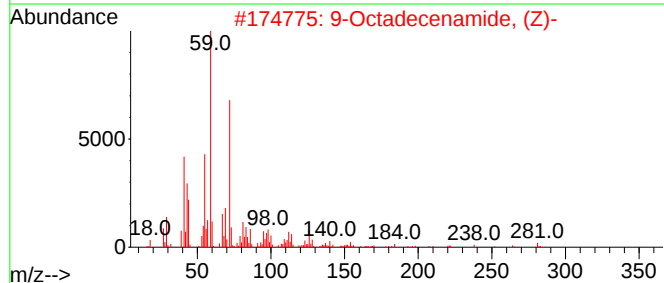
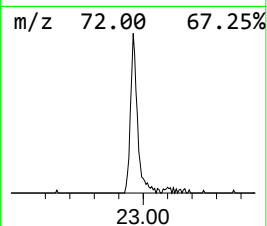
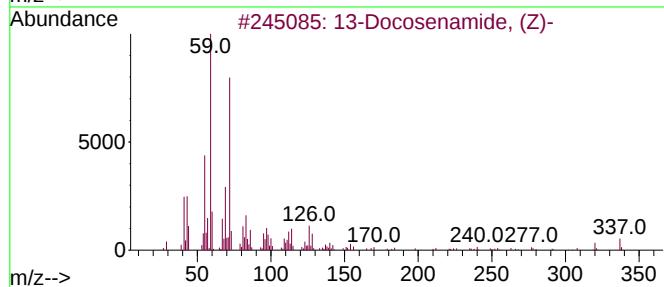
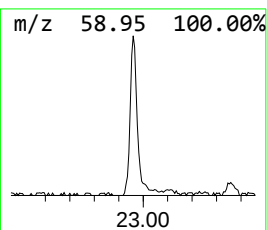
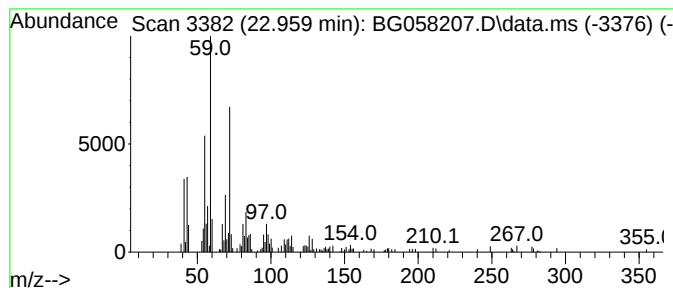
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 13-Docosenamide, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.959	4.40 ng/ul	210927	Chrysene-d12	21.537

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	90
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	83
4			Palmitoleamide	253	C16H31NO	106010-22-4	72
5			Dodecanamide	199	C12H25NO	001120-16-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058207.D
Acq On : 13 Jul 2023 20:52
Operator : CG/JU
Sample : 03414-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A46G0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.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
Cyclohexene	3.158	370.4	ng/u1	5021170	1	7.900	271113	20.0
Tetrachloroethy...	4.621	2.2	ng/u1	29324	1	7.900	271113	20.0
2-Cyclohexen-1-ol	5.802	11.2	ng/u1	151710	1	7.900	271113	20.0
Cyclohexene, 3-...	6.184	3.2	ng/u1	43656	1	7.900	271113	20.0
2-Cyclohexen-1-one	6.525	20.5	ng/u1	277438	1	7.900	271113	20.0
unknown-01	8.076	5.7	ng/u1	77428	1	7.900	271113	20.0
unknown-02	8.153	7.7	ng/u1	103881	1	7.900	271113	20.0
2-Chlorocyclohe...	8.217	56.0	ng/u1	759788	1	7.900	271113	20.0
1,2-Cyclohexane...	8.664	2.1	ng/u1	28968	1	7.900	271113	20.0
Cyclohexane, 1,...	8.899	3.1	ng/u1	41757	1	7.900	271113	20.0
13-Docosenamide...	22.959	4.4	ng/u1	210927	5	21.537	957721	20.0