

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
 Data File : BG058200.D
 Acq On : 13 Jul 2023 15:54
 Operator : CG/JU
 Sample : 03414-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4646

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: BG058200.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	46	rVB	2534672	4974138	100.00%	22.294%
2	3.993	151	154	159	rVB2	7607	11071	0.22%	0.050%
3	4.098	166	172	176	rVB3	4607	7581	0.15%	0.034%
4	4.339	208	213	220	rVB5	11420	19232	0.39%	0.086%
5	4.621	257	261	270	rVB2	16370	26157	0.53%	0.117%
6	4.997	321	325	328	rBV3	4275	6274	0.13%	0.028%
7	5.362	383	387	392	rBV3	9698	16142	0.32%	0.072%
8	5.544	413	418	425	rBV	9625	20691	0.42%	0.093%
9	5.802	453	462	468	rBV3	54787	123599	2.48%	0.554%
10	5.914	476	481	484	rBV4	3529	5017	0.10%	0.022%
11	6.184	521	527	536	rVB2	22744	38832	0.78%	0.174%
12	6.525	579	585	597	rBV	146792	258811	5.20%	1.160%
13	7.071	668	678	690	rBV	218234	396461	7.97%	1.777%
14	7.230	698	705	719	rBV	175519	292735	5.89%	1.312%
15	7.436	731	740	751	rBV	215104	388251	7.81%	1.740%
16	7.518	751	754	761	rVB5	4733	8389	0.17%	0.038%
17	7.618	765	771	777	rBV6	5034	9869	0.20%	0.044%
18	7.906	813	820	827	rBV	141532	239462	4.81%	1.073%
19	7.976	827	832	837	rVV4	9337	14149	0.28%	0.063%
20	8.076	843	849	856	rVV2	26936	49016	0.99%	0.220%
21	8.152	856	862	867	rVV2	39804	69708	1.40%	0.312%
22	8.217	867	873	887	rVV	383618	663314	13.34%	2.973%
23	8.611	932	940	947	rBV	275705	491737	9.89%	2.204%
24	8.728	955	960	969	rVB2	20243	41750	0.84%	0.187%
25	8.899	983	989	995	rBV3	17023	29603	0.60%	0.133%
26	9.063	1010	1017	1029	rBV	207855	375733	7.55%	1.684%
27	9.163	1029	1034	1035	rVV4	4509	7318	0.15%	0.033%
28	9.198	1036	1040	1048	rVB2	8117	16259	0.33%	0.073%
29	9.786	1133	1140	1150	rVB	171622	302299	6.08%	1.355%
30	10.326	1225	1232	1245	rBV	269778	526963	10.59%	2.362%
31	10.708	1287	1297	1304	rBV	219684	399677	8.04%	1.791%
32	10.843	1313	1320	1334	rBV	234840	449193	9.03%	2.013%
33	11.519	1426	1435	1440	rBV8	3192	8265	0.17%	0.037%
34	12.653	1624	1628	1636	rVB5	5619	9259	0.19%	0.041%
35	12.765	1640	1647	1649	rVV3	2731	5324	0.11%	0.024%
36	13.358	1743	1748	1755	rBV2	5048	7446	0.15%	0.033%

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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
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37	13.423	1755	1759	1767	rVV5	4966	7893	0.16%	0.035%
38	13.946	1841	1848	1858	rBV	522877	764539	15.37%	3.427%
39	14.181	1881	1888	1891	rBV4	13736	26380	0.53%	0.118%
40	14.234	1891	1897	1908	rVB	602437	960503	19.31%	4.305%
41	14.539	1942	1949	1959	rBV	368011	585451	11.77%	2.624%
42	14.745	1976	1984	2002	rBV	173657	363220	7.30%	1.628%
43	14.892	2005	2009	2011	rVB5	4038	5283	0.11%	0.024%
44	15.532	2111	2118	2128	rBV	799444	1204294	24.21%	5.398%
45	15.650	2132	2138	2154	rVV	222311	347383	6.98%	1.557%
46	15.767	2154	2158	2163	rVB2	3783	5968	0.12%	0.027%
47	16.078	2206	2211	2216	rBV6	4620	10319	0.21%	0.046%
48	16.972	2359	2363	2369	rBV	62256	85290	1.71%	0.382%
49	17.289	2407	2417	2426	rBV	480990	770232	15.48%	3.452%
50	17.383	2427	2433	2445	rVV	844066	1315619	26.45%	5.896%
51	17.947	2524	2529	2535	rVB3	9068	12629	0.25%	0.057%
52	18.182	2564	2569	2576	rBV7	3369	8918	0.18%	0.040%
53	18.717	2657	2660	2663	rBV3	3480	5774	0.12%	0.026%
54	18.746	2663	2665	2671	rVB3	3844	6172	0.12%	0.028%
55	18.969	2699	2703	2708	rBV5	7255	11997	0.24%	0.054%
56	19.116	2725	2728	2731	rBB3	8906	8118	0.16%	0.036%
57	19.239	2744	2749	2755	rBV4	15220	26564	0.53%	0.119%
58	19.674	2817	2823	2834	rBV	1061511	1557192	31.31%	6.979%
59	20.144	2899	2903	2905	rBV4	4041	5061	0.10%	0.023%
60	20.567	2971	2975	2978	rBV6	4940	8473	0.17%	0.038%
61	20.990	3044	3047	3053	rBV6	17773	23130	0.47%	0.104%
62	21.267	3092	3094	3100	rVB2	8510	11205	0.23%	0.050%
63	21.425	3117	3121	3126	rBV2	19405	23320	0.47%	0.105%
64	21.537	3134	3140	3150	rBV	487374	857237	17.23%	3.842%
65	21.719	3168	3171	3175	rVB4	13654	17541	0.35%	0.079%
66	22.312	3269	3272	3277	rVB5	14662	20101	0.40%	0.090%
67	22.965	3378	3383	3395	rVB5	74734	173661	3.49%	0.778%
68	23.764	3514	3519	3525	rVB2	23096	43819	0.88%	0.196%
69	24.469	3627	3639	3653	rBV2	587321	1679482	33.76%	7.527%
70	24.686	3665	3676	3688	rVB	336850	976291	19.63%	4.376%
71	25.785	3856	3863	3870	rVB4	24422	62362	1.25%	0.280%
72	27.077	4081	4083	4085	rBV3	9983	10832	0.22%	0.049%

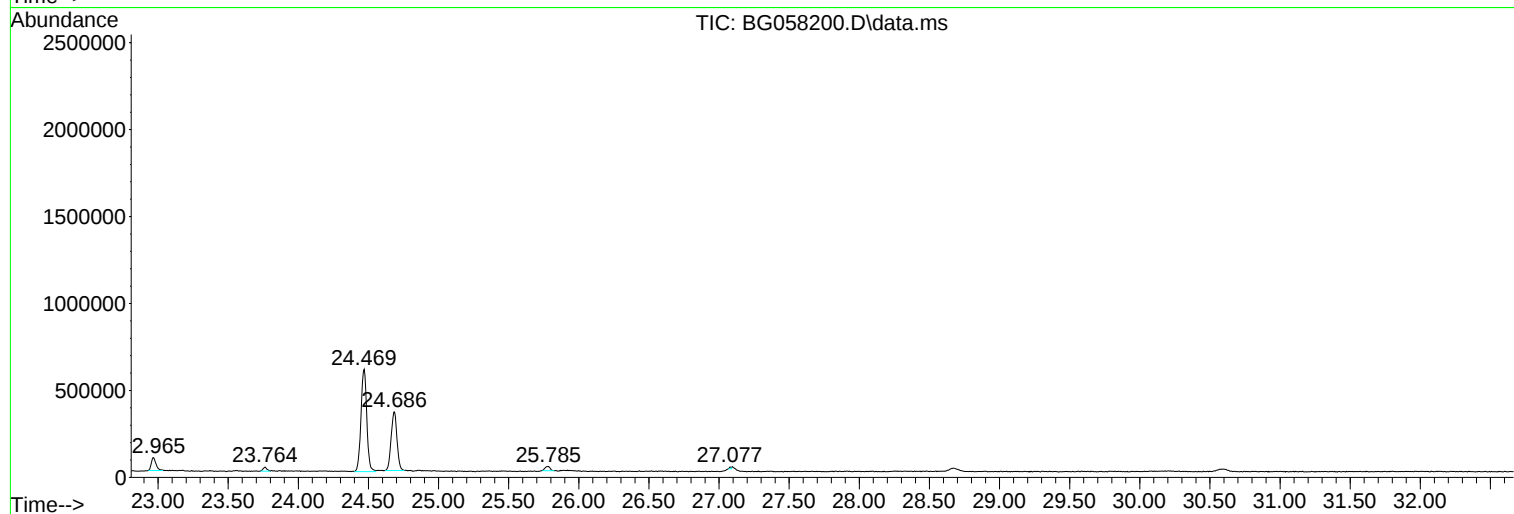
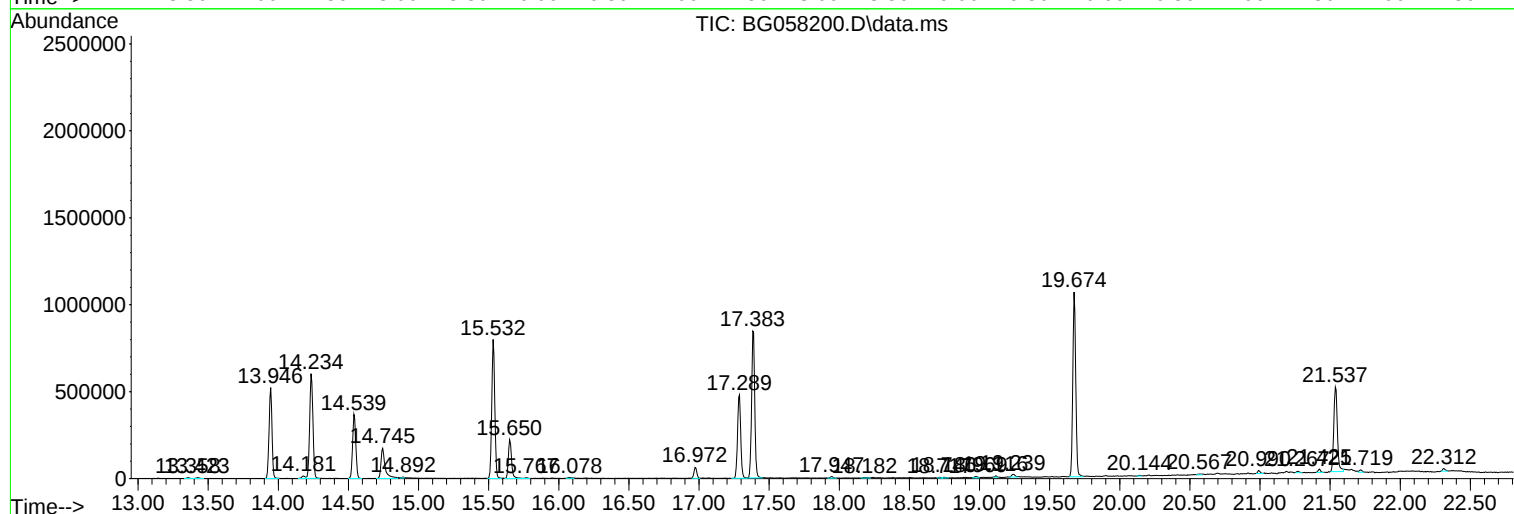
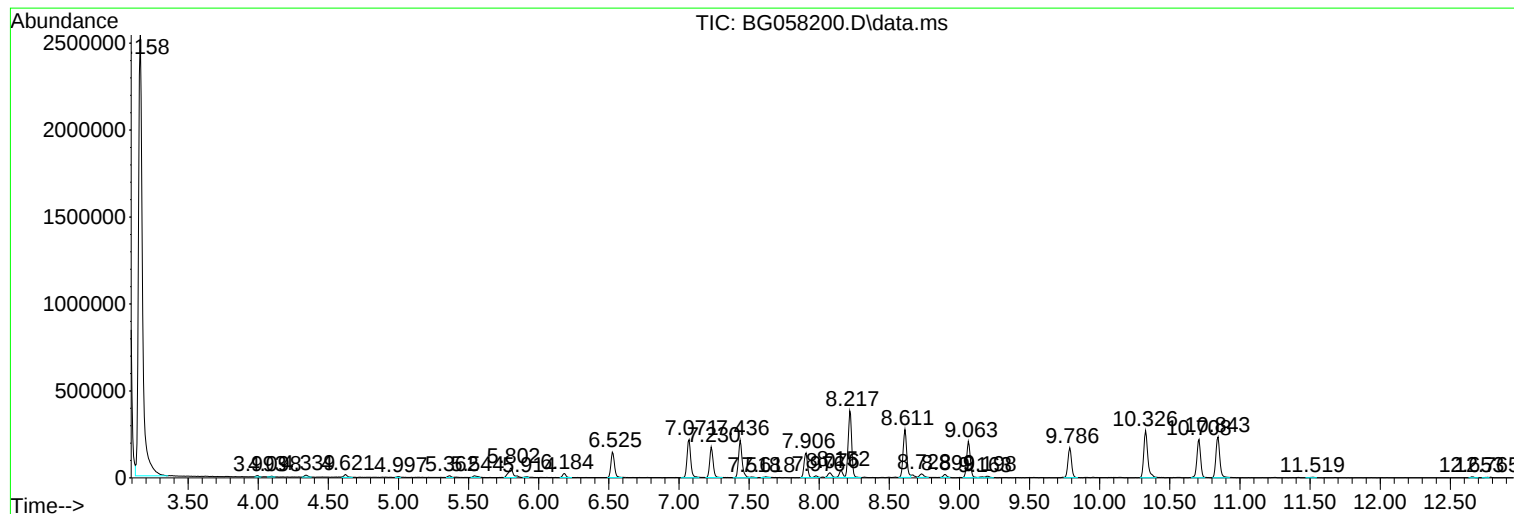
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : CG/JU
Sample : 03414-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4646

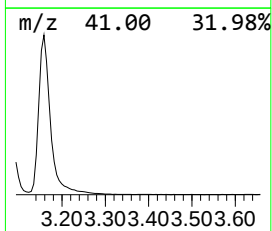
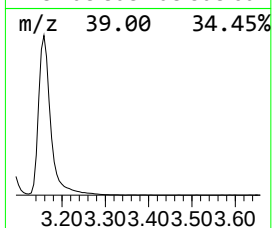
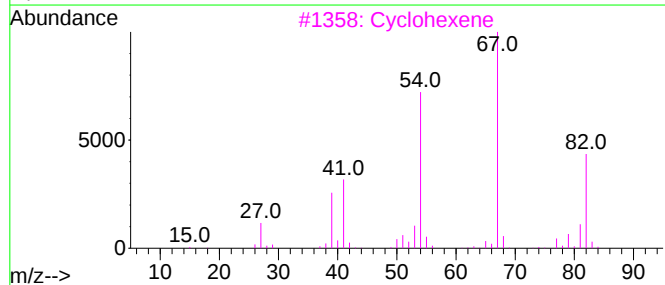
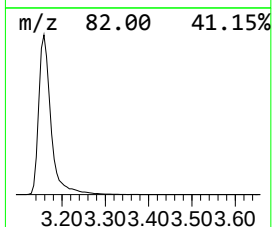
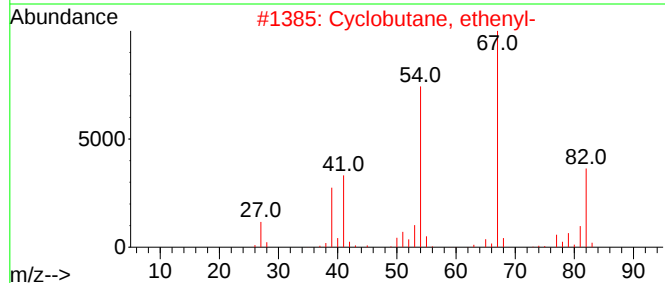
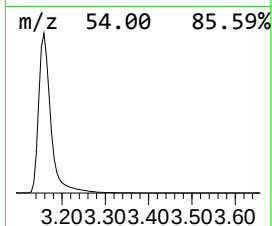
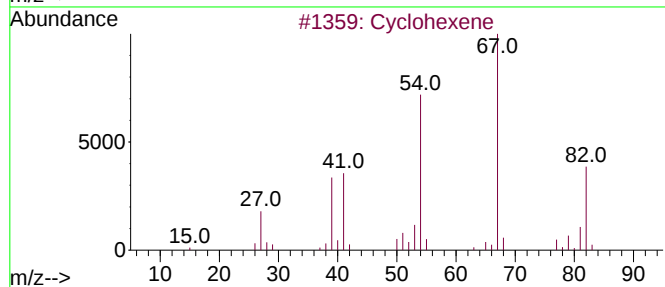
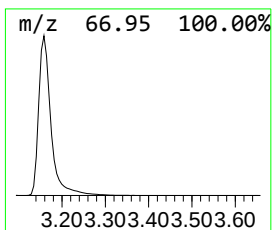
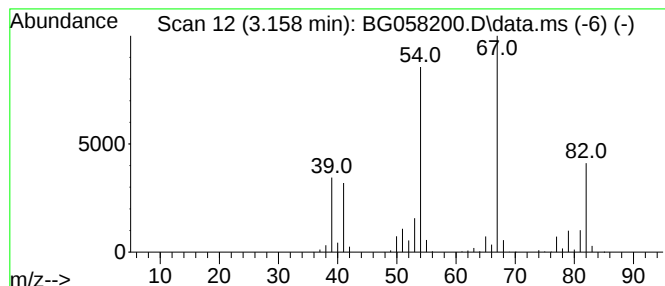
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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	415.44 ng/ul	4974140	1,4-Dichlorobenzene-d4	7.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene	82	C6H10	000110-83-8	95
2			Cyclobutane, ethenyl-	82	C6H10	002597-49-1	94
3			Cyclohexene	82	C6H10	000110-83-8	91
4			Cyclohexene	82	C6H10	000110-83-8	91
5			Ethylidenecyclobutane	82	C6H10	001528-21-8	91



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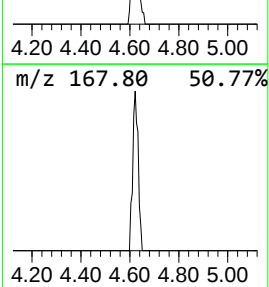
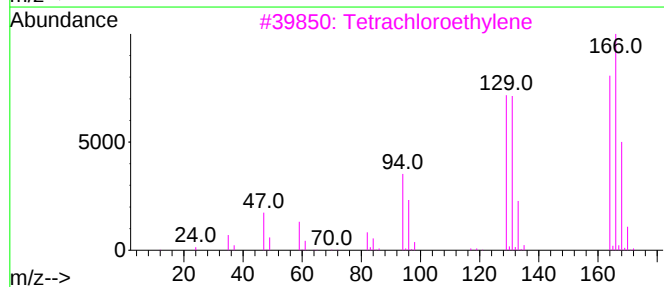
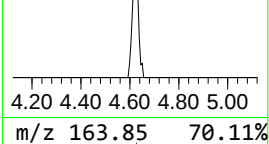
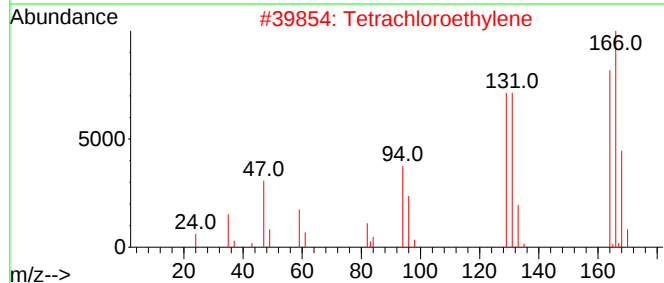
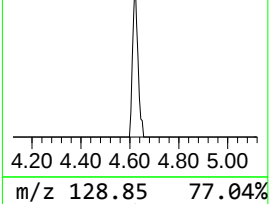
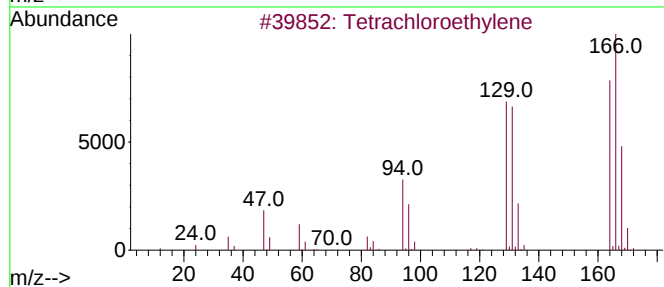
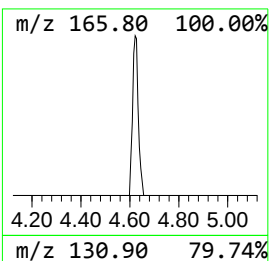
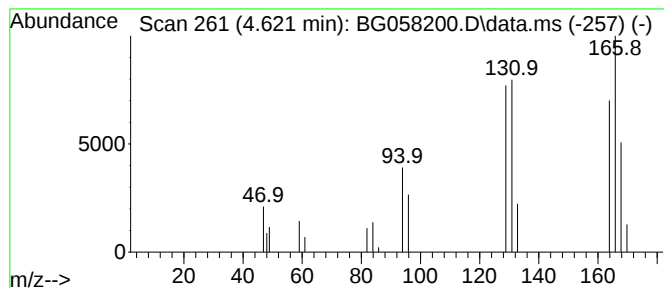
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Tetrachloroethylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.621	2.18 ng/ul	26157	1,4-Dichlorobenzene-d4	7.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	43



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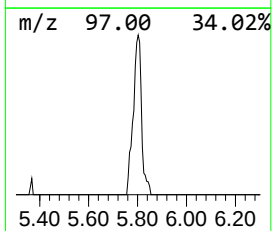
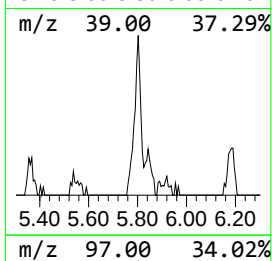
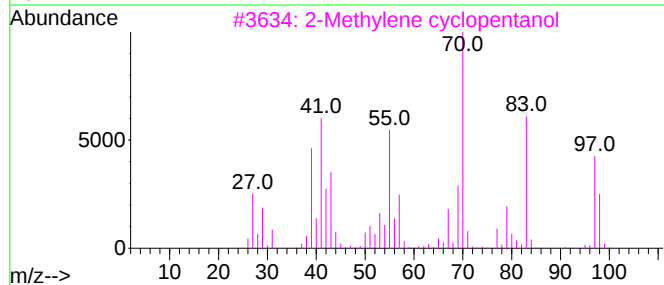
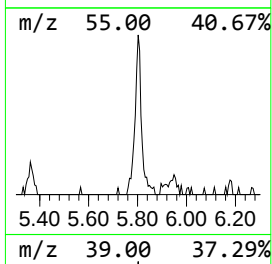
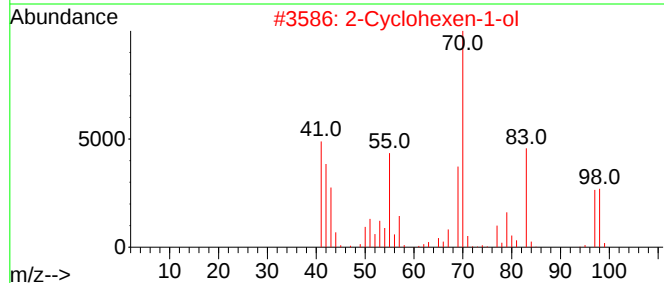
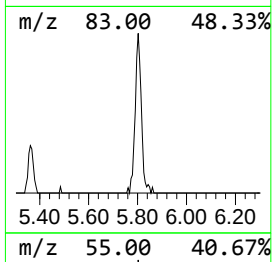
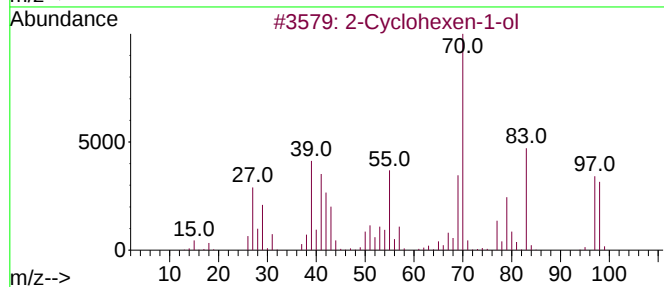
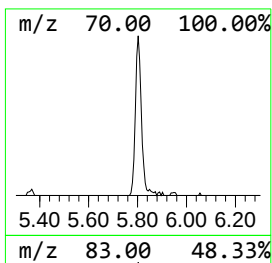
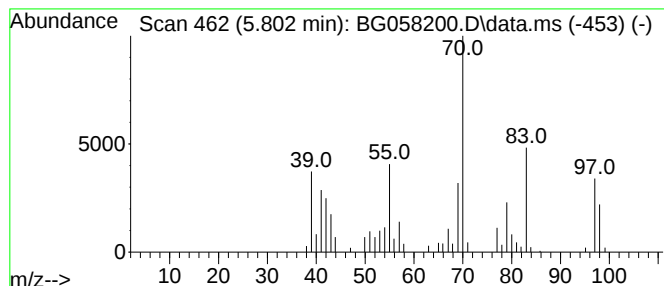
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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	10.32 ng/ul	123599	1,4-Dichlorobenzene-d4	7.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	96
2	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	87
3	2-Methylene cyclopentanol			98	C6H10O	020461-31-8	80
4	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	76
5	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	38



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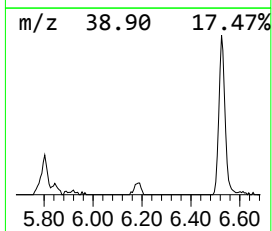
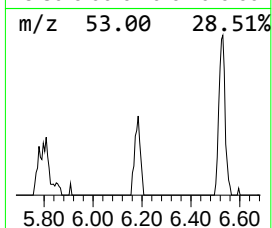
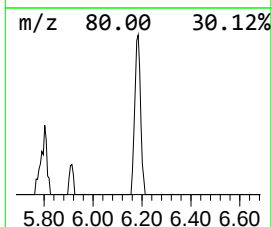
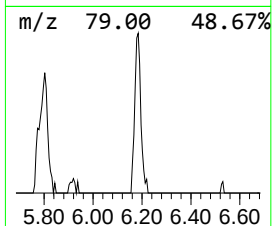
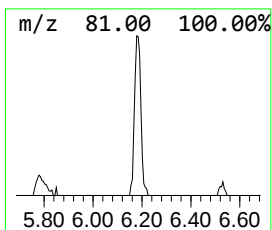
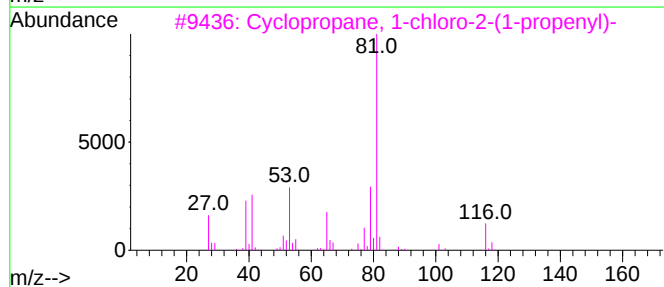
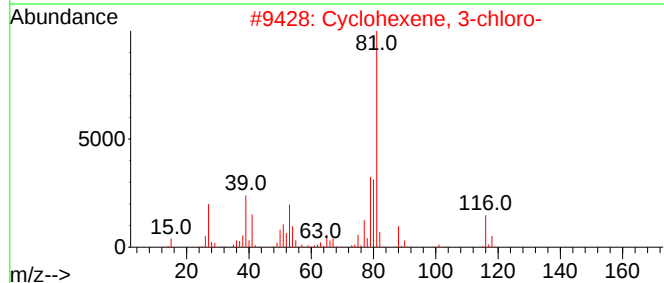
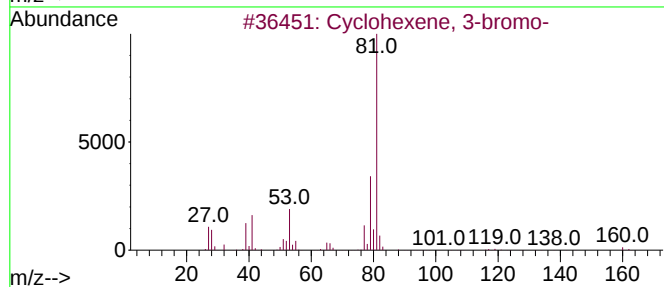
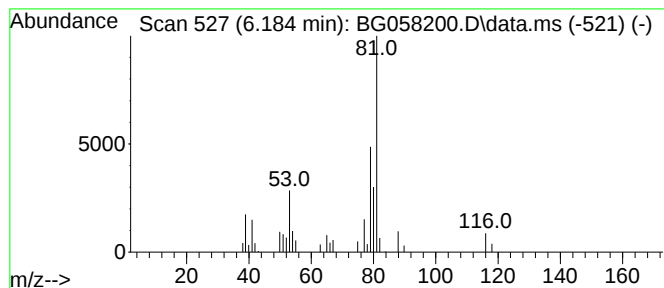
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.184	3.24 ng/ul	38832	1,4-Dichlorobenzene-d4	7.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 3-bromo-	160	C6H9Br	001521-51-3	72
2			Cyclohexene, 3-chloro-	116	C6H9Cl	002441-97-6	64
3			Cyclopropane, 1-chloro-2-(1-prop...	116	C6H9Cl	074752-94-6	64
4			1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	64
5			Cyclopropane, 1-chloro-2-ethenyl...	116	C6H9Cl	062337-93-3	64



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Data File : BG058200.D
Acq On : 13 Jul 2023 15:54
Operator : CG/JU
Sample : 03414-03
Misc :
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Instrument :
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ClientSampleId :
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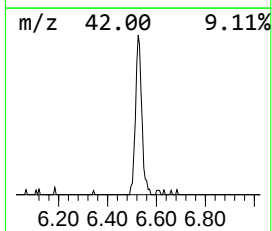
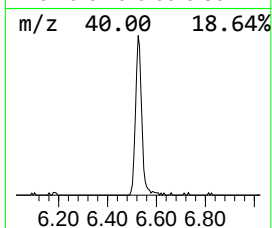
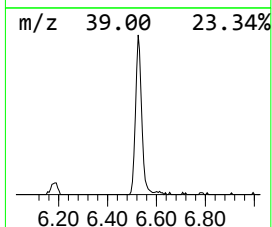
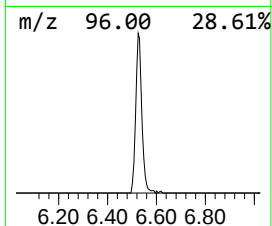
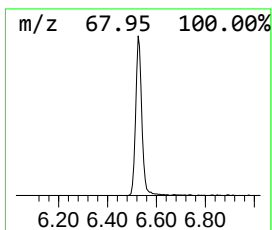
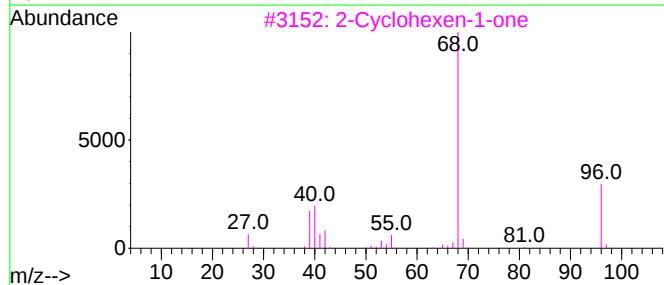
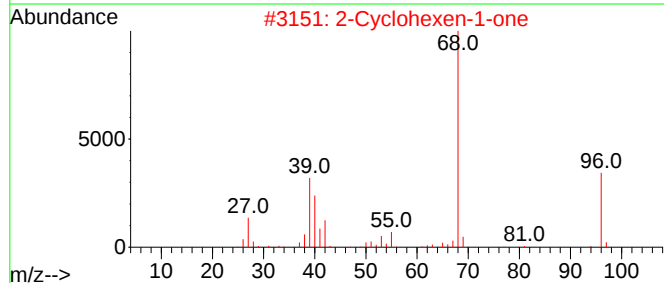
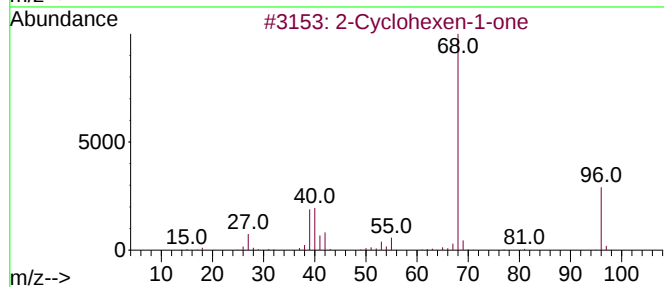
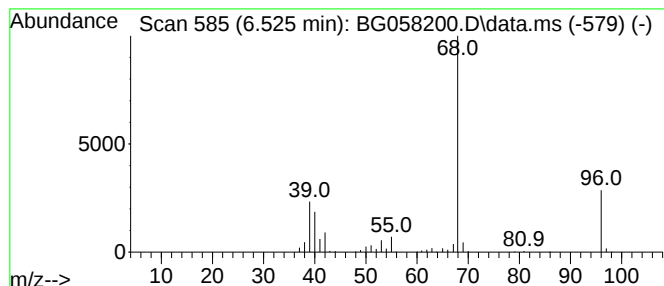
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.525	21.62 ng/ul	258811	1,4-Dichlorobenzene-d4	7.906

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	91
5			2H-Pyran-2-one, 5,6-dihydro-	98	C5H6O2	003393-45-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058200.D
Acq On : 13 Jul 2023 15:54
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Sample : 03414-03
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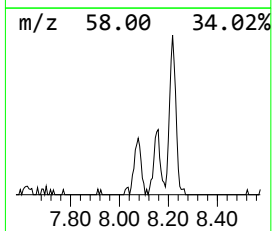
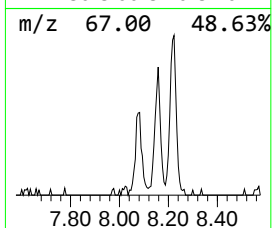
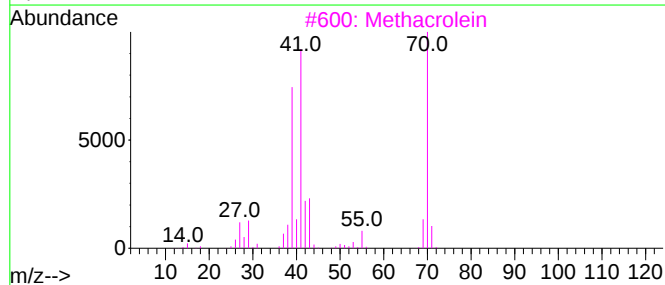
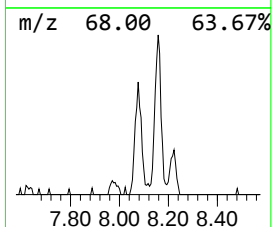
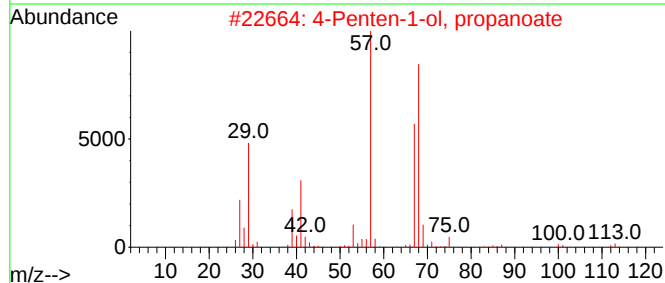
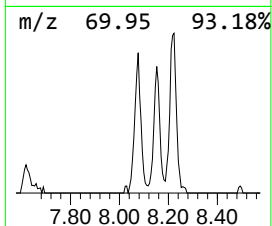
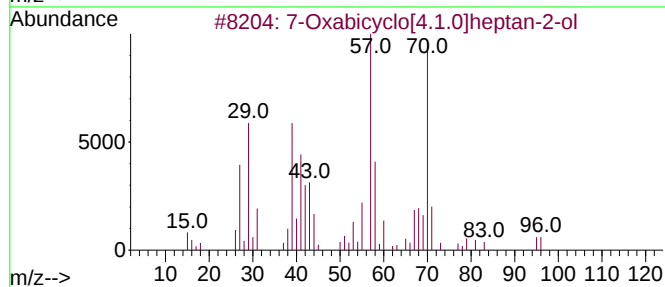
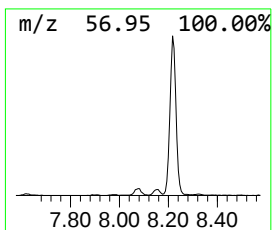
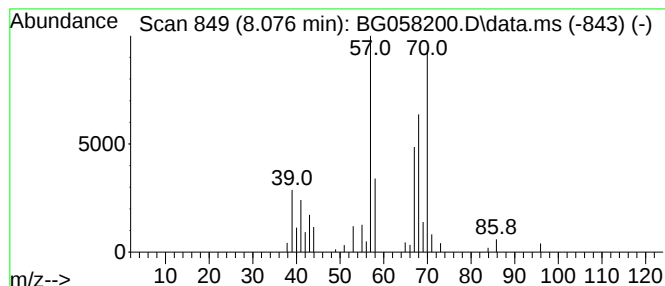
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Quant Title : SVOA CALIBRATION

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	4.09 ng/ul	49016	1,4-Dichlorobenzene-d4	7.906

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		cis,cis-2,7-Nonadiene	124	C9H16	036901-84-5	25
5		1,4-Pentanediamine	102	C5H14N2	000591-77-5	23



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058200.D
Acq On : 13 Jul 2023 15:54
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Sample : 03414-03
Misc :
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Instrument :
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ClientSampleId :
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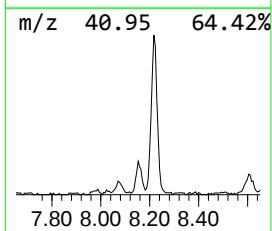
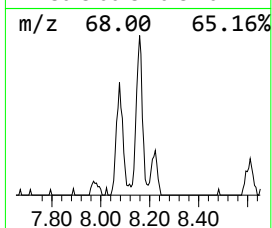
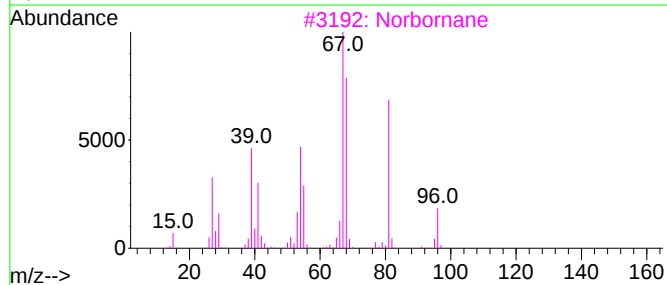
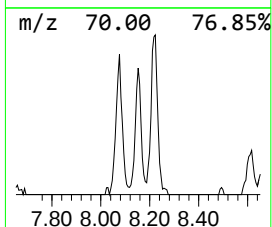
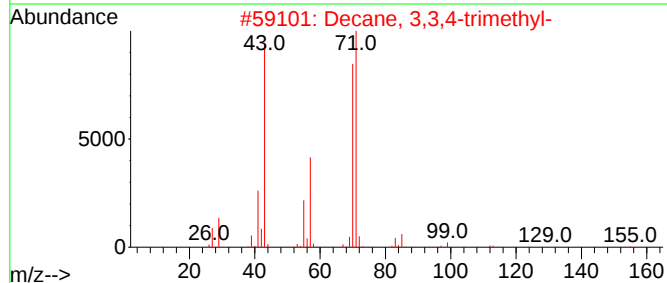
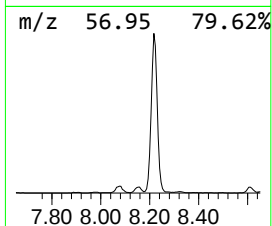
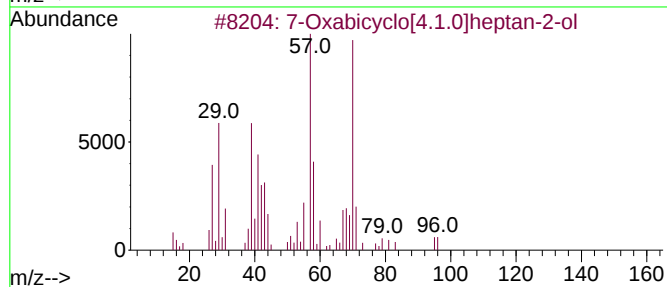
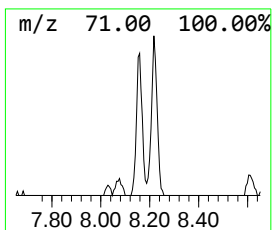
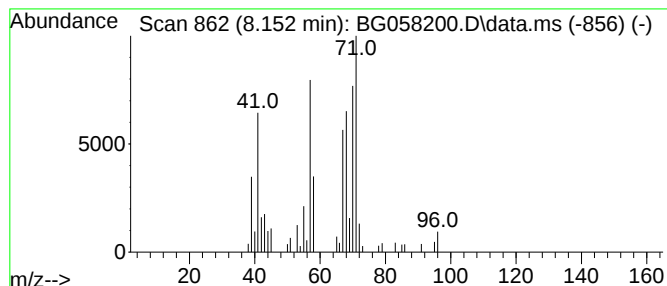
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.152	5.82 ng/ul	69708	1,4-Dichlorobenzene-d4	7.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Oxabicyclo[4.1.0]heptan-2-ol	114	C6H10O2	001192-78-5	38
2			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	25
3			Norbornane	96	C7H12	000279-23-2	14
4			2-Propen-1-amine, N-ethyl-	85	C5H11N	002424-02-4	12
5			Butanoic acid, 4-pentenyl ester	156	C9H16O2	030563-31-6	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058200.D
Acq On : 13 Jul 2023 15:54
Operator : CG/JU
Sample : 03414-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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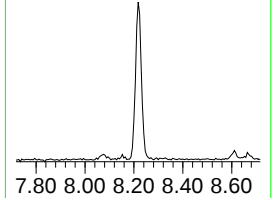
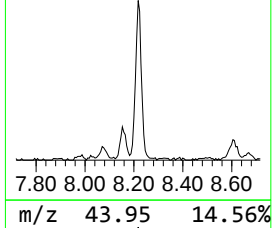
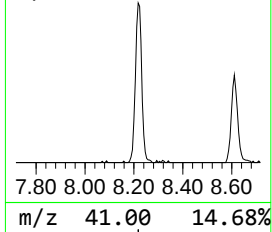
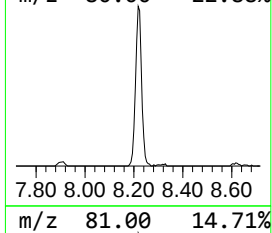
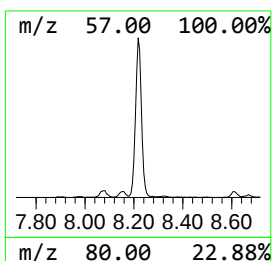
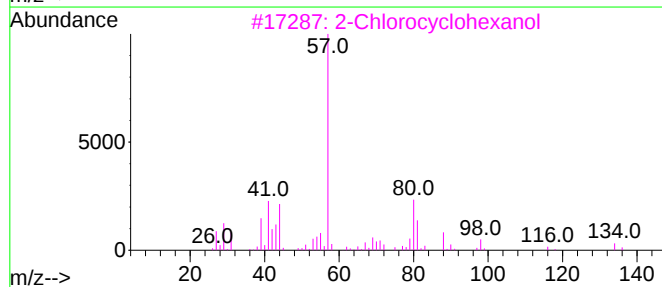
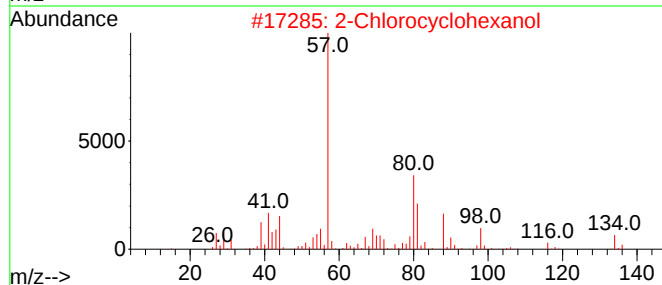
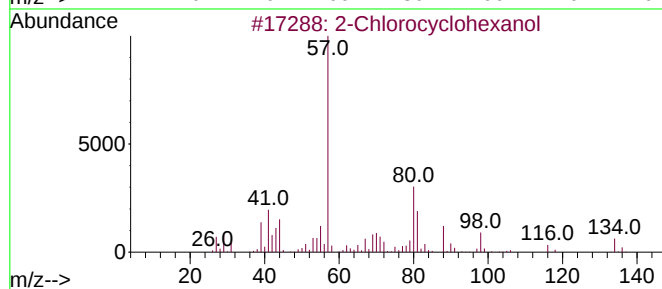
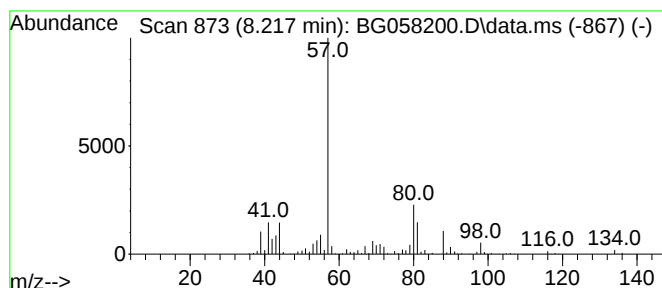
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Chlorocyclohexanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.217	55.40 ng/ul	663314	1,4-Dichlorobenzene-d4	7.906

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	94	
2	2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	94	
3	2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	83	
4	2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	53	
5	2-Chlorocyclohexanol	134	C6H11ClO	001561-86-0	43	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058200.D
Acq On : 13 Jul 2023 15:54
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Sample : 03414-03
Misc :
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Instrument :
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ClientSampleId :
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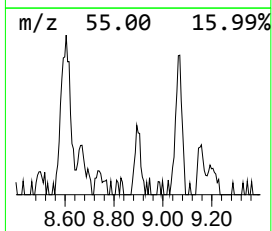
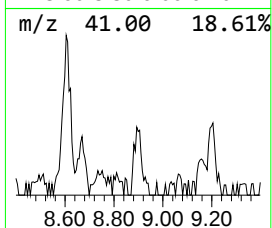
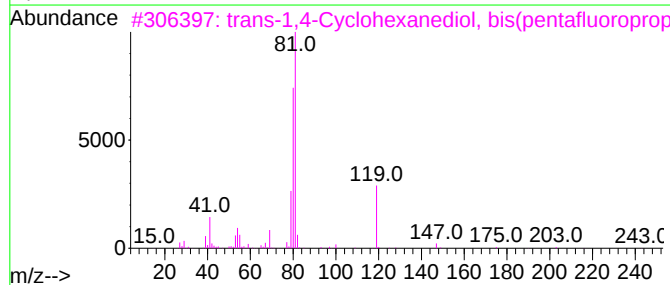
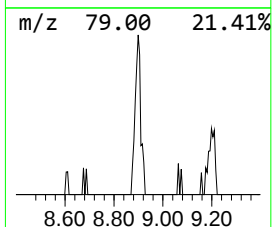
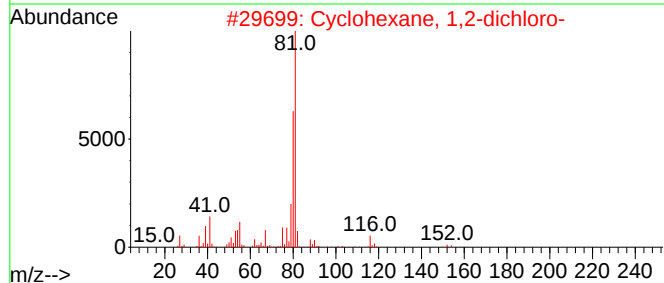
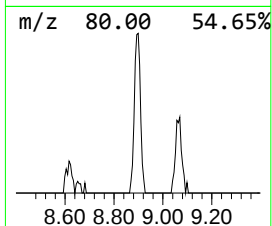
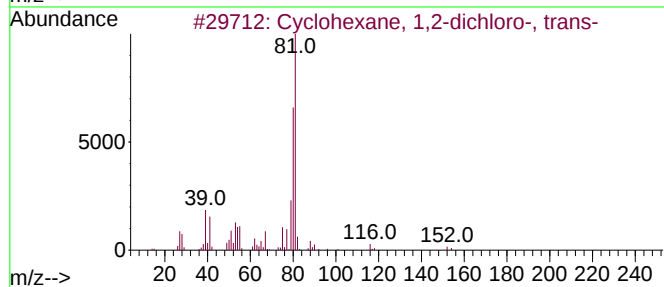
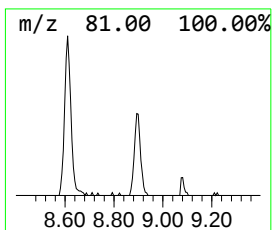
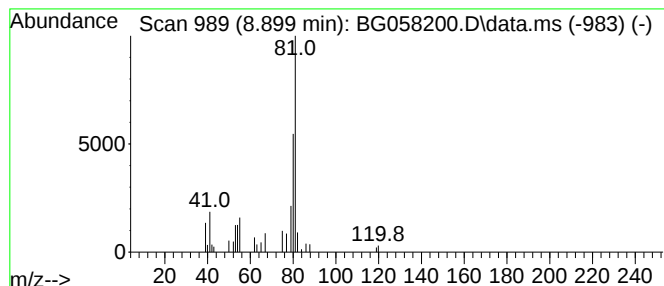
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.899	2.47 ng/ul	29603	1,4-Dichlorobenzene-d4	7.906

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1,2-dichloro-, trans-	152	C6H10Cl2	000822-86-6	83
2		Cyclohexane, 1,2-dichloro-	152	C6H10Cl2	001121-21-7	72
3		trans-1,4-Cyclohexanediol, bis(p...	408	C12H10F10O4	1000384-30-6	72
4		Bi-2-cyclohexen-1-yl	162	C12H18	001541-20-4	64
5		Cyclohexane, 1,3-dichloro-, cis-	152	C6H10Cl2	024955-63-3	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058200.D
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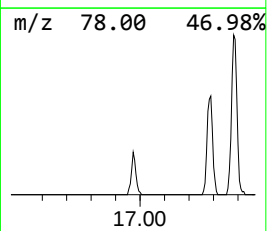
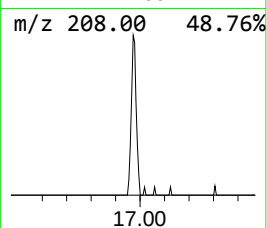
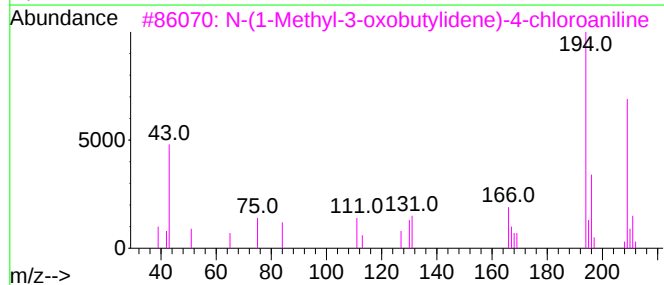
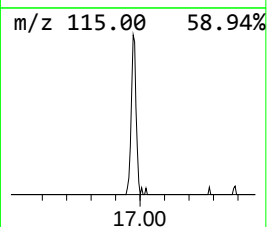
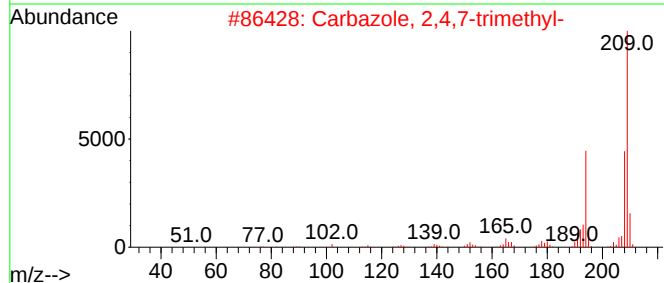
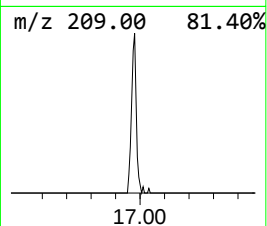
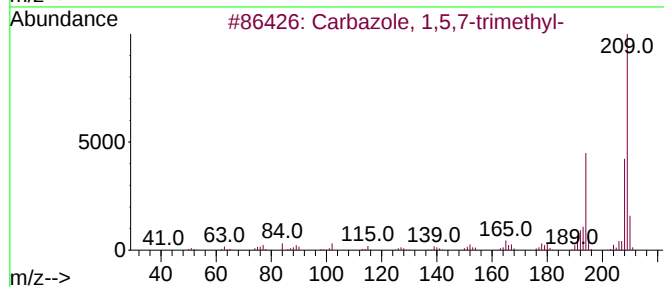
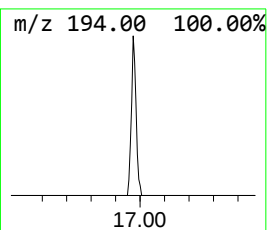
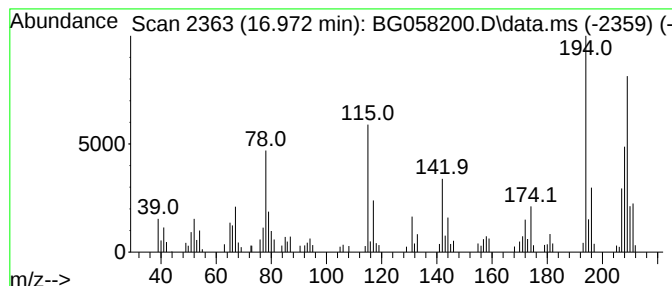
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Carbazole, 1,5,7-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.971	2.21 ng/ul	85290	Phenanthrene-d10	17.289

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbazole, 1,5,7-trimethyl-	209	C15H15N	078787-84-5	52
2			Carbazole, 2,4,7-trimethyl-	209	C15H15N	078787-89-0	49
3			N-(1-Methyl-3-oxobutylidene)-4-c...	209	C11H12ClNO	050519-24-9	49
4			2-Amino-4-trimethylsilylbenzoic ...	209	C10H15NO2Si	030127-88-9	43
5			4-(Dimethylamino)phenol, TMS	209	C11H19NOSi	1000504-21-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
Data File : BG058200.D
Acq On : 13 Jul 2023 15:54
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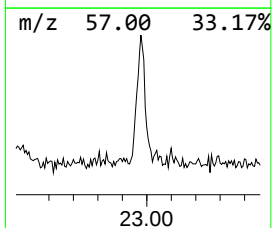
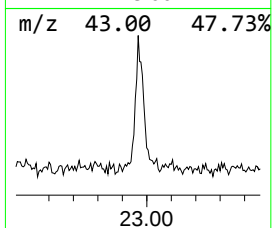
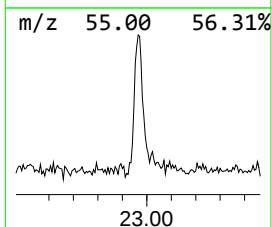
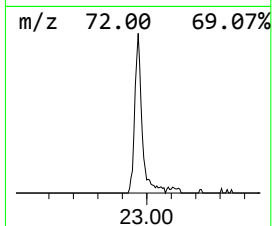
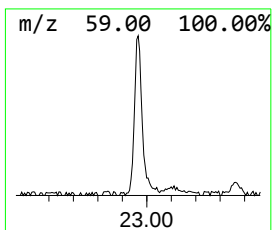
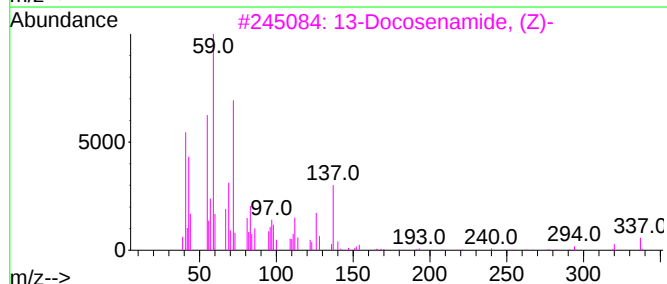
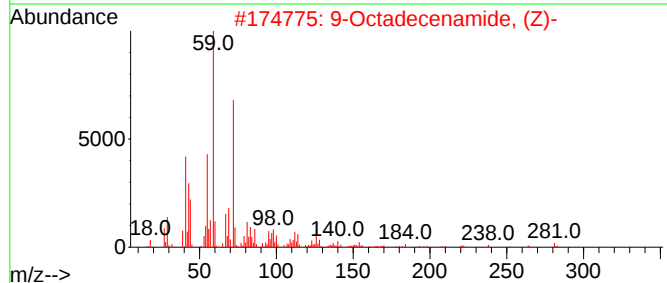
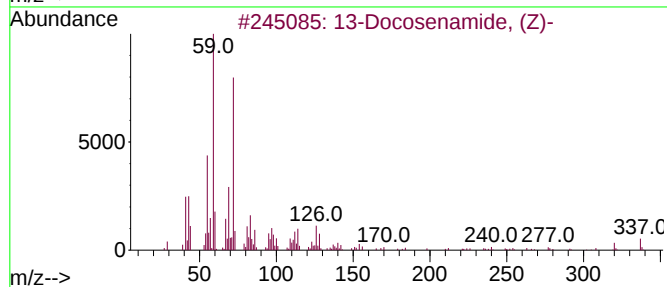
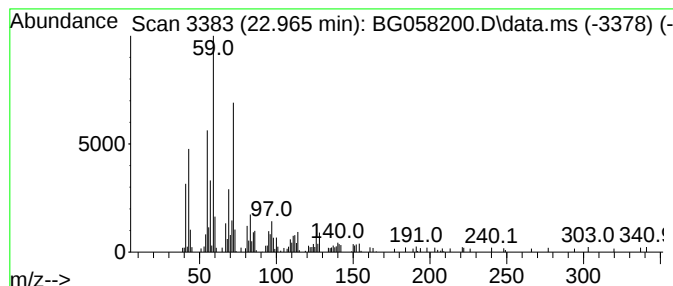
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.965	4.05 ng/ul	173661	Chrysene-d12	21.537

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	94
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	86
3		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	81
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
5		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071323\
 Data File : BG058200.D
 Acq On : 13 Jul 2023 15:54
 Operator : CG/JU
 Sample : 03414-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4646

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	415.4	ng/ul	4974140	1	7.906	239462	20.0
Tetrachloroethy...	4.621	2.2	ng/ul	26157	1	7.906	239462	20.0
2-Cyclohexen-1-ol	5.802	10.3	ng/ul	123599	1	7.906	239462	20.0
Cyclohexene, 3-...	6.184	3.2	ng/ul	38832	1	7.906	239462	20.0
2-Cyclohexen-1-one	6.525	21.6	ng/ul	258811	1	7.906	239462	20.0
unknown-01	8.076	4.1	ng/ul	49016	1	7.906	239462	20.0
unknown-02	8.152	5.8	ng/ul	69708	1	7.906	239462	20.0
2-Chlorocyclohe...	8.217	55.4	ng/ul	663314	1	7.906	239462	20.0
Cyclohexane, 1,...	8.899	2.5	ng/ul	29603	1	7.906	239462	20.0
Carbazole, 1,5,...	16.971	2.2	ng/ul	85290	4	17.289	770232	20.0
13-Docosenamide...	22.965	4.0	ng/ul	173661	5	21.537	857237	20.0