

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
 Data File : BG058187.D
 Acq On : 12 Jul 2023 15:11
 Operator : CG/JU
 Sample : 03387-11
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EX7Z5

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: BG058187.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.141	4	9	16	rBV	948633	1271610	39.08%	7.176%
2	3.217	16	22	51	rVB	1866240	3253855	100.00%	18.362%
3	4.052	160	164	170	rVB3	9119	14071	0.43%	0.079%
4	4.158	178	182	187	rVB3	4793	6486	0.20%	0.037%
5	4.398	219	223	232	rVB6	9580	15553	0.48%	0.088%
6	4.692	267	273	278	rVB4	12193	20673	0.64%	0.117%
7	5.062	332	336	340	rBV4	4570	6808	0.21%	0.038%
8	5.432	394	399	404	rBV5	5222	8066	0.25%	0.046%
9	5.609	423	429	442	rBV3	9564	22383	0.69%	0.126%
10	5.867	465	473	478	rBV3	60159	119800	3.68%	0.676%
11	5.979	489	492	498	rBV5	4217	4660	0.14%	0.026%
12	6.249	532	538	544	rVB2	22904	35602	1.09%	0.201%
13	6.590	589	596	606	rBV	146533	256375	7.88%	1.447%
14	7.131	682	688	702	rVB	138544	244650	7.52%	1.381%
15	7.295	710	716	724	rBV	107473	169694	5.22%	0.958%
16	7.501	744	751	761	rBV	131639	236013	7.25%	1.332%
17	7.589	763	766	769	rVB5	3194	3819	0.12%	0.022%
18	7.689	776	783	786	rBV7	4019	7187	0.22%	0.041%
19	7.971	823	831	838	rBV	163533	271018	8.33%	1.529%
20	8.047	839	844	848	rBV3	8237	12773	0.39%	0.072%
21	8.141	855	860	866	rVB2	27845	49518	1.52%	0.279%
22	8.223	867	874	878	rBV2	33618	56175	1.73%	0.317%
23	8.288	878	885	893	rVB	359912	613661	18.86%	3.463%
24	8.394	899	903	909	rVB5	2517	4189	0.13%	0.024%
25	8.564	928	932	936	rVB6	2767	4210	0.13%	0.024%
26	8.676	943	951	957	rVV	169190	304074	9.35%	1.716%
27	8.729	957	960	966	rVV4	11273	20804	0.64%	0.117%
28	8.793	966	971	981	rVB	23257	44006	1.35%	0.248%
29	8.964	994	1000	1007	rVB	18674	32355	0.99%	0.183%
30	9.134	1022	1029	1037	rVV	131829	235304	7.23%	1.328%
31	9.222	1039	1044	1045	rVV3	4055	4408	0.14%	0.025%
32	9.269	1048	1052	1057	rVB5	8424	14181	0.44%	0.080%
33	9.857	1145	1152	1160	rVB	107552	190180	5.84%	1.073%
34	9.962	1166	1170	1176	rBV3	1768	3678	0.11%	0.021%
35	10.027	1176	1181	1188	rBV8	3141	6844	0.21%	0.039%
36	10.397	1237	1244	1259	rBV	173800	330408	10.15%	1.865%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
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37	10.773	1296	1308	1316	rBV	245112	453741	13.94%	2.561%
38	10.908	1323	1331	1342	rBV	146334	271414	8.34%	1.532%
39	11.584	1437	1446	1452	rBV6	2570	7803	0.24%	0.044%
40	12.718	1635	1639	1645	rVB4	4836	8132	0.25%	0.046%
41	12.824	1653	1657	1659	rBV3	3096	4845	0.15%	0.027%
42	12.853	1659	1662	1666	rVB2	2825	3977	0.12%	0.022%
43	13.411	1753	1757	1763	rBV3	4181	6026	0.19%	0.034%
44	13.482	1763	1769	1773	rBV3	5178	8153	0.25%	0.046%
45	13.999	1851	1857	1863	rBV	330447	465615	14.31%	2.628%
46	14.246	1893	1899	1900	rBV3	5723	8568	0.26%	0.048%
47	14.293	1900	1907	1916	rVB	379476	607357	18.67%	3.427%
48	14.598	1951	1959	1967	rBV	426817	660950	20.31%	3.730%
49	14.798	1986	1993	2011	rBV	95566	201738	6.20%	1.138%
50	14.951	2015	2019	2022	rVV5	4226	7078	0.22%	0.040%
51	15.591	2121	2128	2135	rBV	479567	741213	22.78%	4.183%
52	15.703	2141	2147	2157	rBV	118607	180087	5.53%	1.016%
53	16.179	2224	2228	2234	rVV7	2155	4893	0.15%	0.028%
54	16.872	2339	2346	2351	rVB6	2309	4054	0.12%	0.023%
55	17.025	2367	2372	2377	rBV	45585	65122	2.00%	0.367%
56	17.125	2384	2389	2393	rVB6	1698	3367	0.10%	0.019%
57	17.342	2419	2426	2435	rBV	555542	829655	25.50%	4.682%
58	17.436	2435	2442	2450	rBV	507672	782072	24.04%	4.413%
59	17.630	2473	2475	2482	rVB5	3397	4241	0.13%	0.024%
60	17.994	2533	2537	2542	rVB2	8823	10699	0.33%	0.060%
61	18.223	2571	2576	2579	rBV4	3084	5149	0.16%	0.029%
62	18.558	2629	2633	2637	rVB4	2555	4112	0.13%	0.023%
63	18.770	2665	2669	2670	rVV2	3175	3776	0.12%	0.021%
64	18.799	2672	2674	2678	rVB3	3987	5072	0.16%	0.029%
65	18.952	2695	2700	2701	rBV4	2524	3462	0.11%	0.020%
66	19.022	2708	2712	2716	rVB3	6606	9077	0.28%	0.051%
67	19.075	2716	2721	2724	rBV7	2670	4911	0.15%	0.028%
68	19.163	2732	2736	2739	rBV4	8304	8633	0.27%	0.049%
69	19.293	2754	2758	2762	rBV6	10157	14838	0.46%	0.084%
70	19.634	2811	2816	2823	rBV9	3415	8009	0.25%	0.045%
71	19.728	2823	2832	2841	rBV	633045	930253	28.59%	5.250%
72	19.880	2854	2858	2862	rBV5	2299	3720	0.11%	0.021%
73	19.945	2867	2869	2875	rVV3	3318	4225	0.13%	0.024%
74	20.021	2877	2882	2884	rBV6	2625	4074	0.13%	0.023%
75	20.192	2908	2911	2914	rVB5	3123	3968	0.12%	0.022%
76	20.603	2979	2981	2984	rBV4	3381	3992	0.12%	0.023%

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77	20.673	2989	2993	2995	rBV3	3815	4955	0.15%	0.028%
78	20.903	3026	3032	3033	rBV5	2763	4240	0.13%	0.024%
79	20.961	3040	3042	3048	rVB6	4781	7100	0.22%	0.040%
80	21.038	3051	3055	3060	rVV6	10637	14056	0.43%	0.079%
81	21.473	3126	3129	3135	rVB	17510	23821	0.73%	0.134%
82	21.596	3143	3150	3159	rBV2	527442	861018	26.46%	4.859%
83	21.713	3166	3170	3175	rBV8	4693	6781	0.21%	0.038%
84	22.377	3279	3283	3288	rVB5	9243	14206	0.44%	0.080%
85	23.041	3389	3396	3411	rBV3	75246	173893	5.34%	0.981%
86	23.846	3528	3533	3543	rVB2	22127	46247	1.42%	0.261%
87	24.563	3644	3655	3668	rBV2	327098	960540	29.52%	5.420%
88	24.792	3682	3694	3708	rBV	331348	1006753	30.94%	5.681%
89	25.897	3873	3882	3892	rVB3	32166	94452	2.90%	0.533%
90	26.202	3930	3934	3936	rBV5	3392	5458	0.17%	0.031%
91	27.236	4100	4110	4125	rVB2	32328	111705	3.43%	0.630%
92	28.846	4374	4384	4397	rVB5	22373	89563	2.75%	0.505%
93	30.779	4705	4713	4717	rBV8	14149	42571	1.31%	0.240%

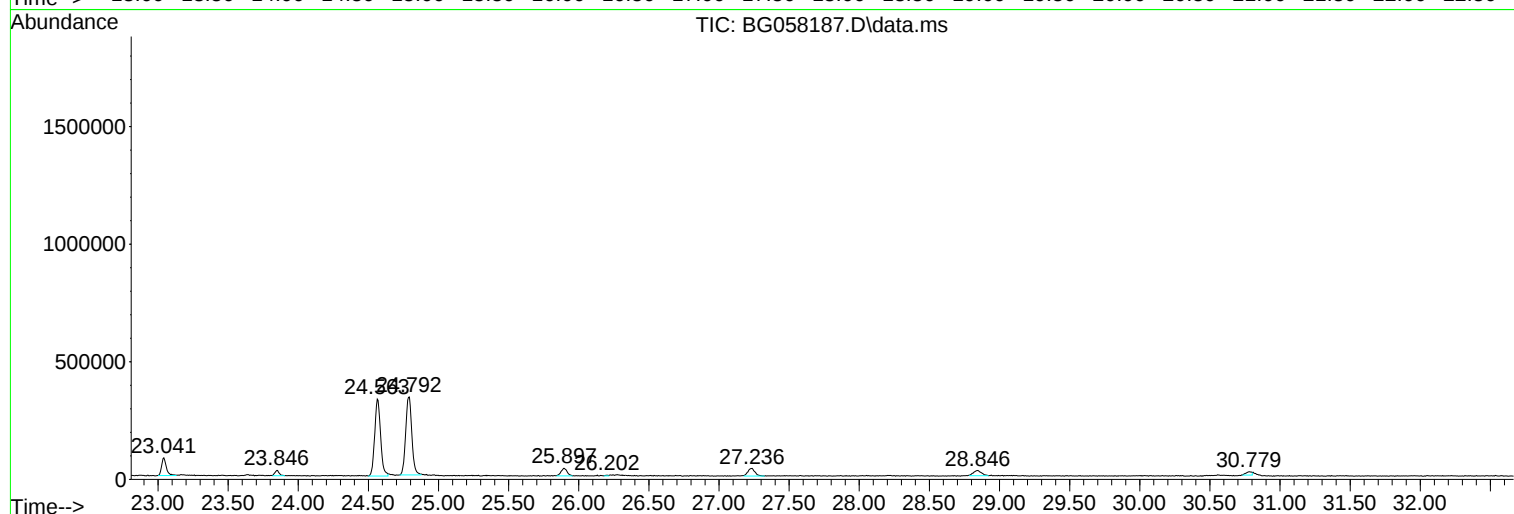
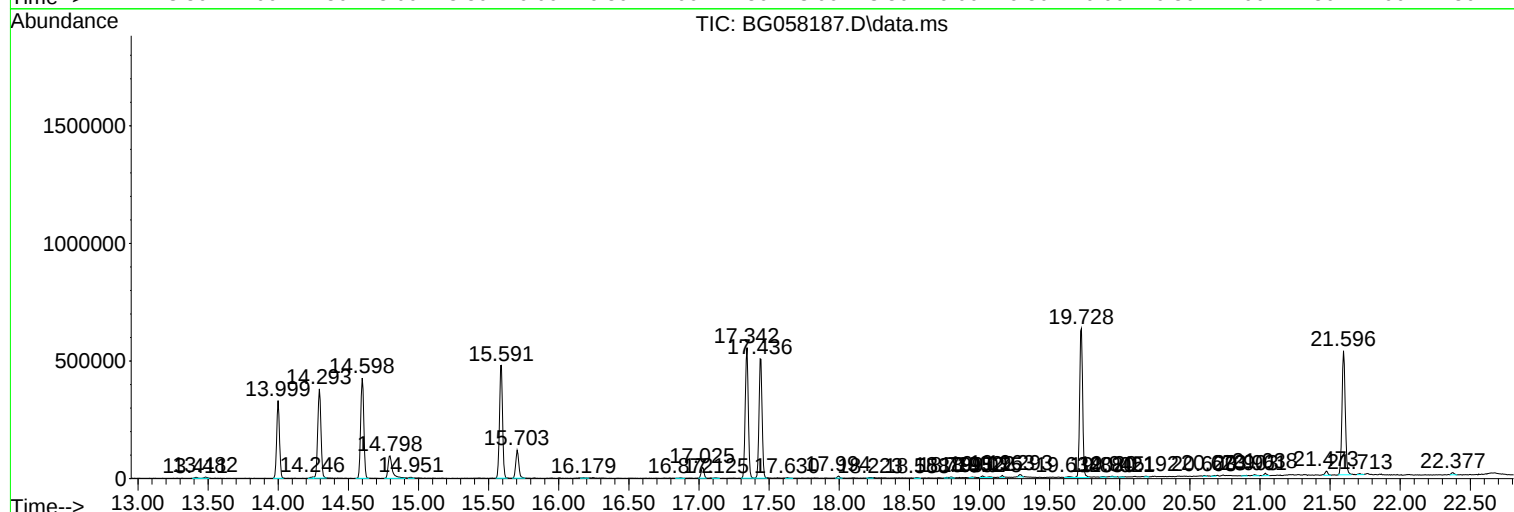
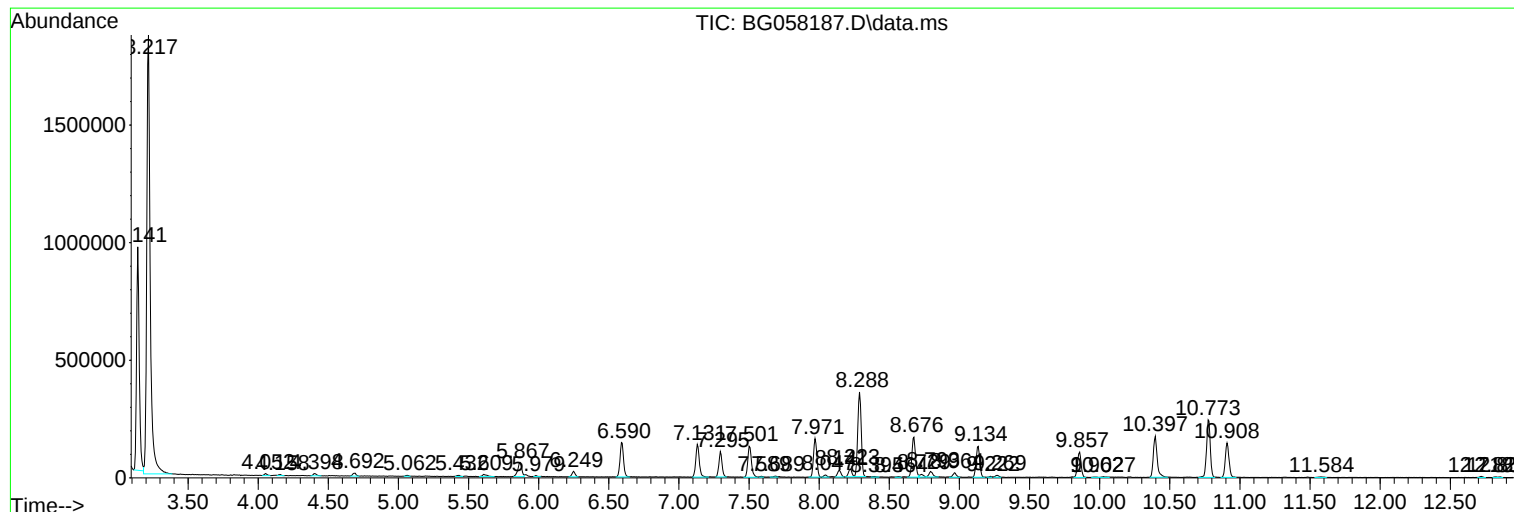
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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



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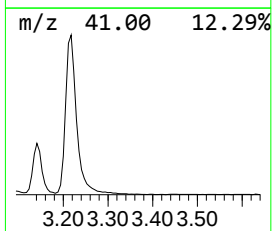
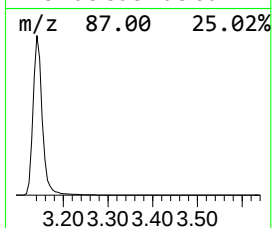
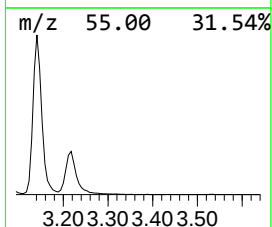
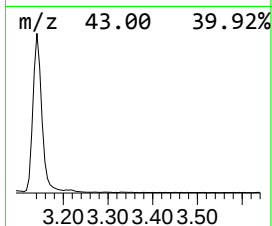
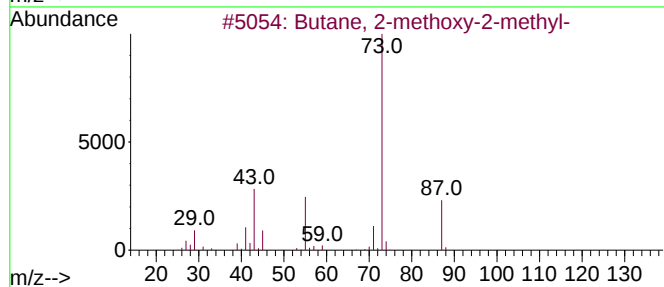
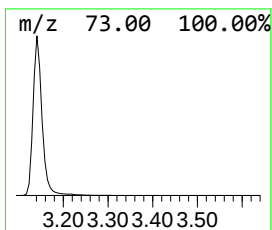
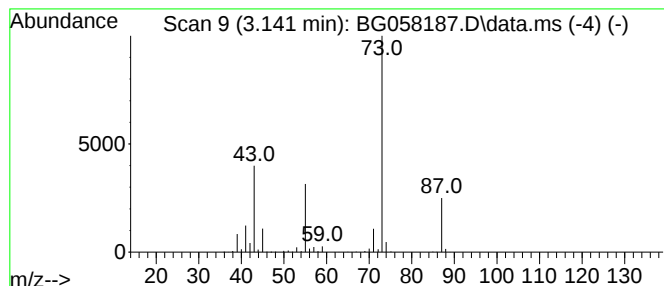
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.141	93.84 ng/ul	1271610	1,4-Dichlorobenzene-d4	7.971

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27



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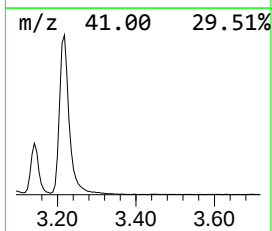
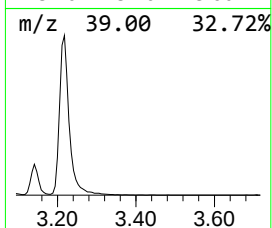
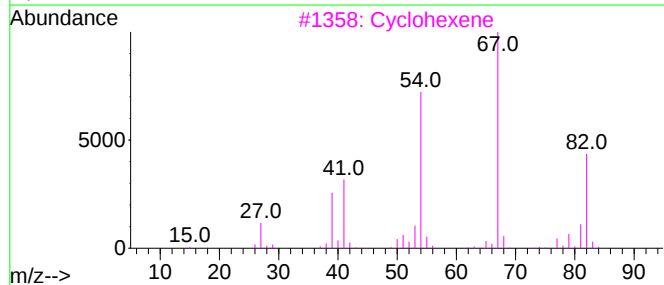
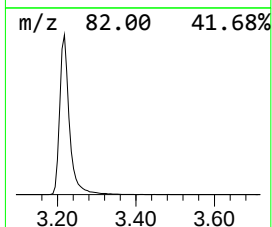
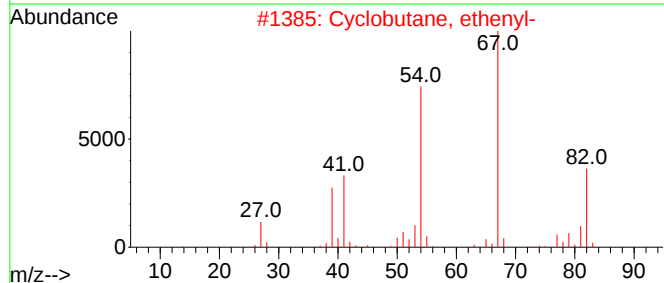
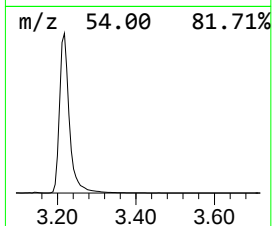
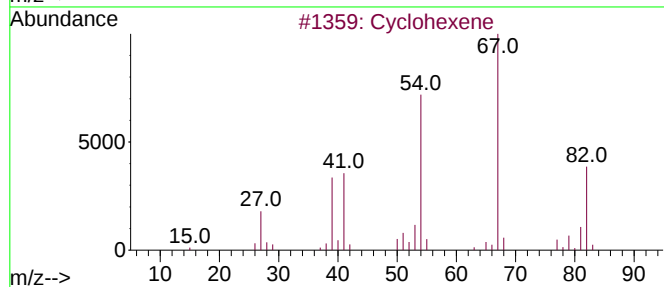
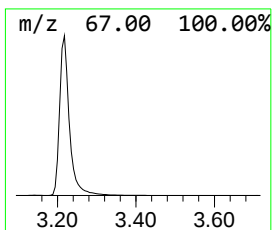
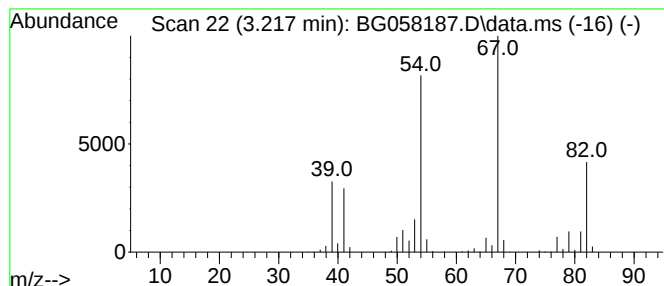
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclohexene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.217	240.12 ng/ul	3253860	1,4-Dichlorobenzene-d4	7.971

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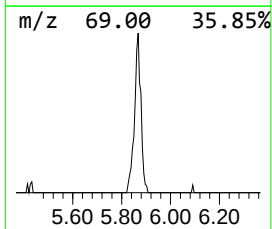
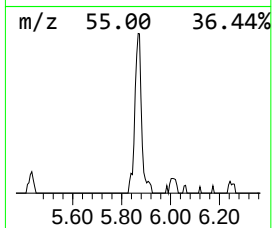
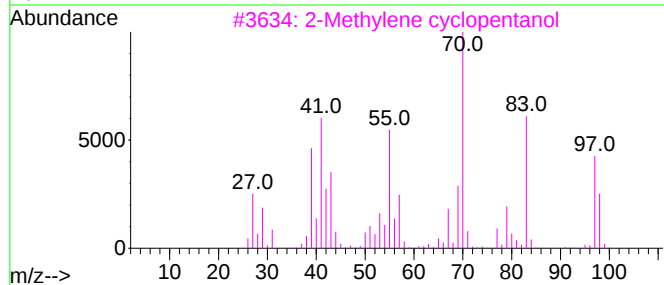
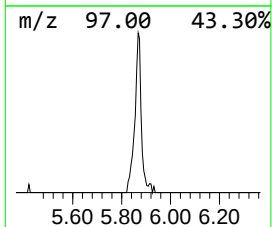
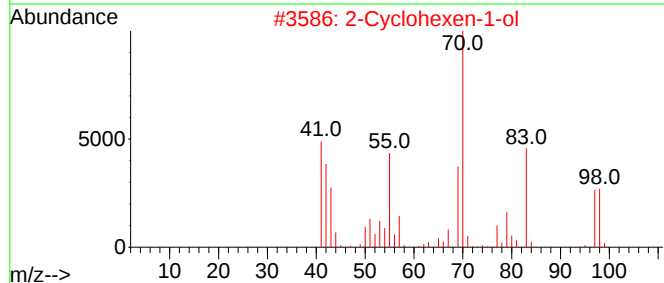
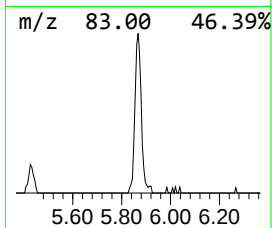
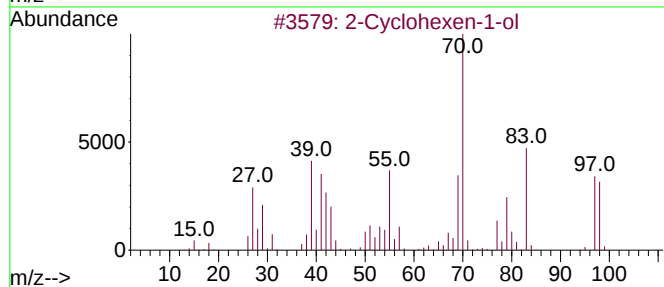
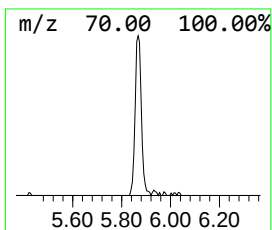
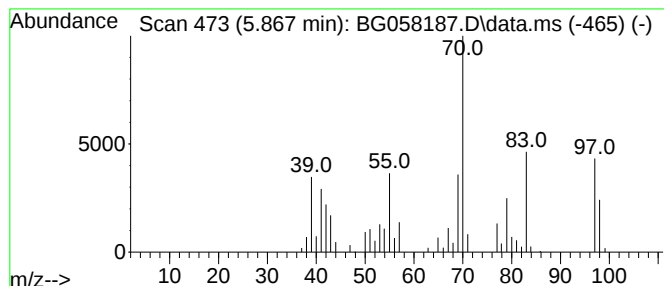
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.867	8.84 ng/ul	119800	1,4-Dichlorobenzene-d4	7.971

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	94
2	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	81
3	2-Methylene cyclopentanol			98	C6H10O	020461-31-8	80
4	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	70
5	2-Cyclohexen-1-ol			98	C6H10O	000822-67-3	58



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ClientSampleId :
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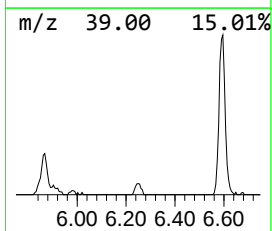
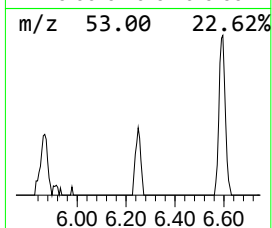
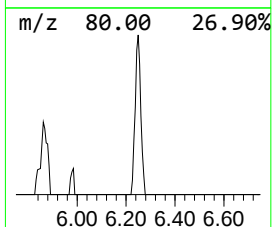
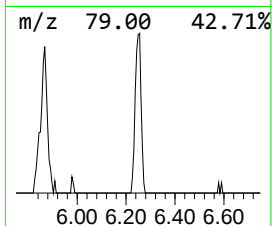
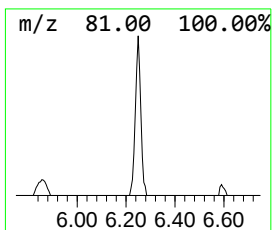
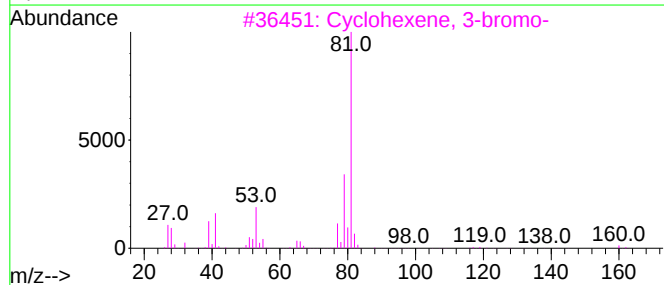
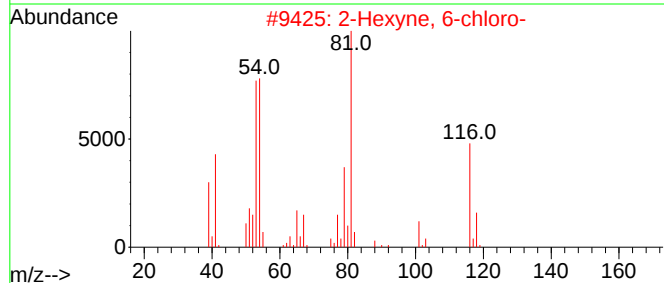
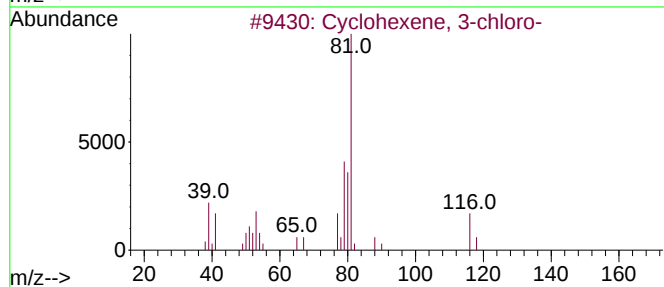
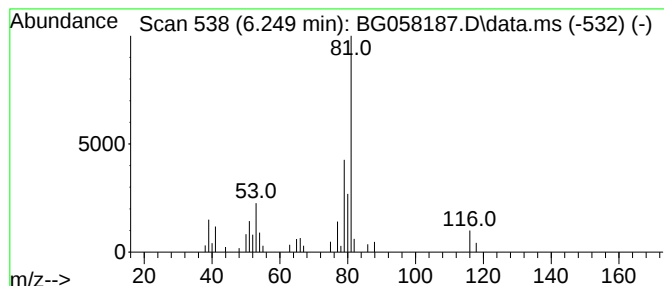
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclohexene, 3-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.249	2.63 ng/ul	35602	1,4-Dichlorobenzene-d4	7.971

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 3-chloro-	116	C6H9Cl	002441-97-6	91
2			2-Hexyne, 6-chloro-	116	C6H9Cl	028077-73-8	72
3			Cyclohexene, 3-bromo-	160	C6H9Br	001521-51-3	64
4			3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	59
5			Cyclopropane, 1-chloro-2-ethenyl...	116	C6H9Cl	062337-93-3	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
Data File : BG058187.D
Acq On : 12 Jul 2023 15:11
Operator : CG/JU
Sample : 03387-11
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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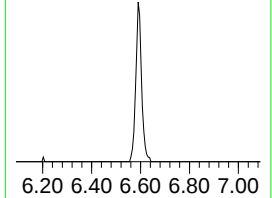
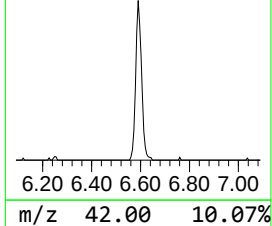
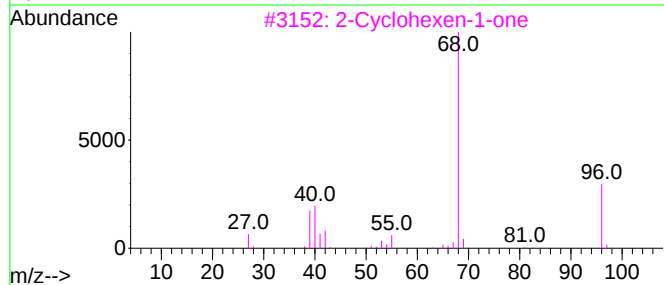
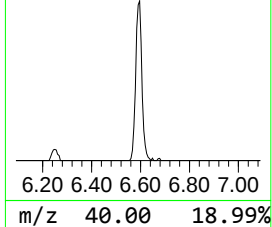
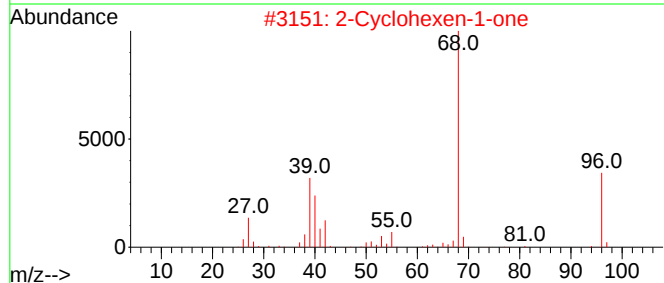
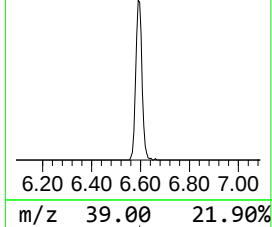
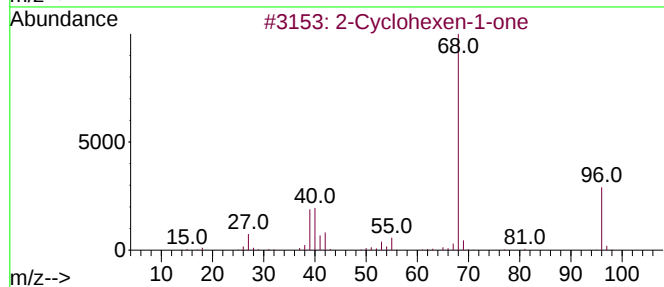
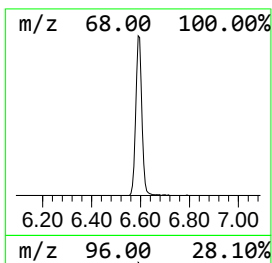
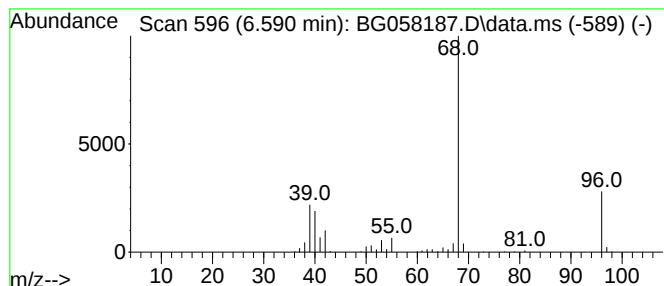
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Cyclohexen-1-one Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.590	18.92 ng/ul	256375	1,4-Dichlorobenzene-d4	7.971

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\BG071123\
Data File : BG058187.D
Acq On : 12 Jul 2023 15:11
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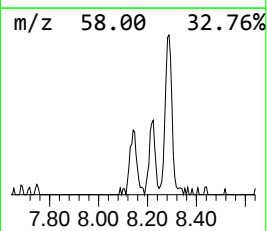
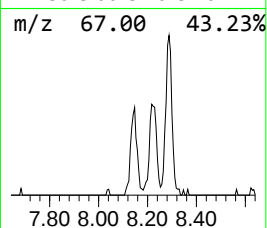
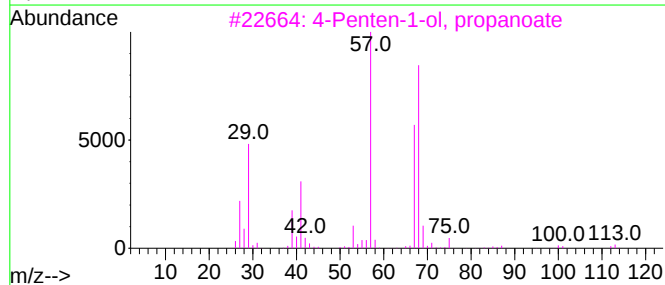
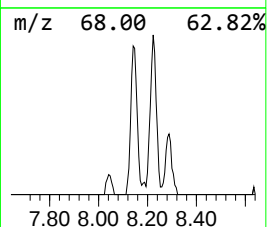
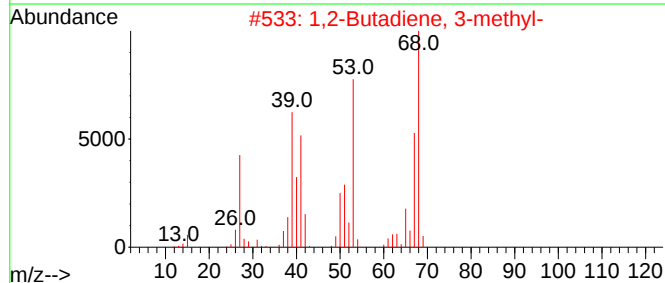
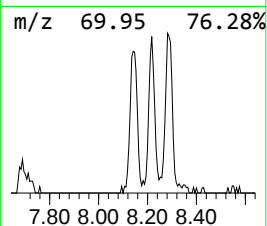
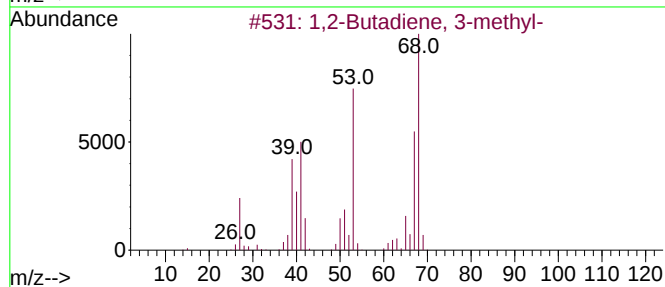
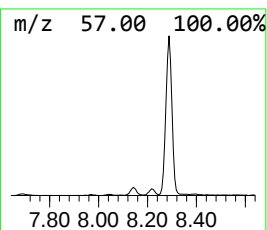
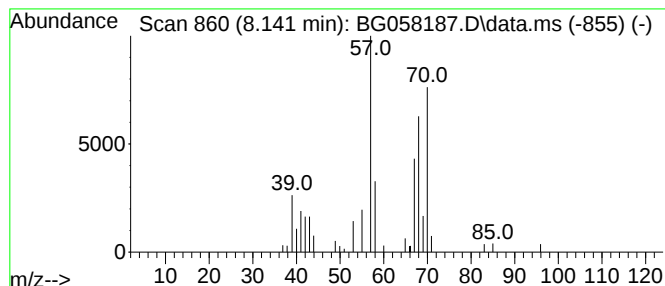
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.141	3.65 ng/ul	49518	1,4-Dichlorobenzene-d4	7.971

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Butadiene, 3-methyl-	68	C5H8	000598-25-4	27
2			1,2-Butadiene, 3-methyl-	68	C5H8	000598-25-4	22
3			4-Penten-1-ol, propanoate	142	C8H14O2	030563-30-5	17
4			cis,cis-2,7-Nonadiene	124	C9H16	036901-84-5	16
5			2,3-Octadiene, 7-methyl-	124	C9H16	061129-33-7	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
Data File : BG058187.D
Acq On : 12 Jul 2023 15:11
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Instrument :
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ClientSampleId :
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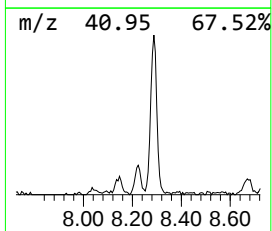
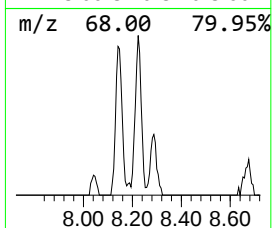
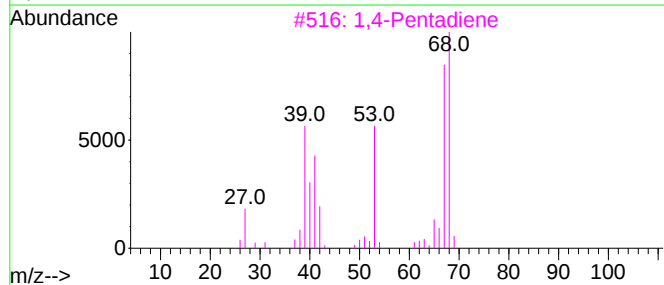
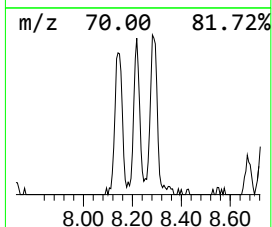
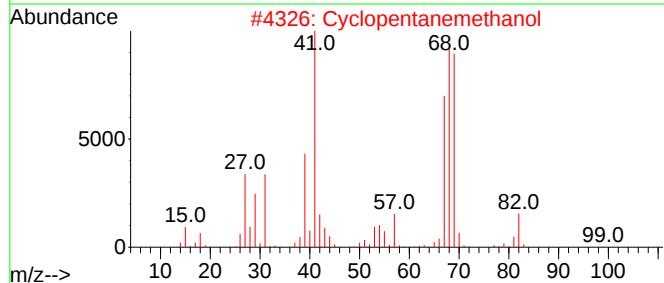
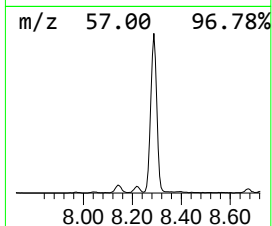
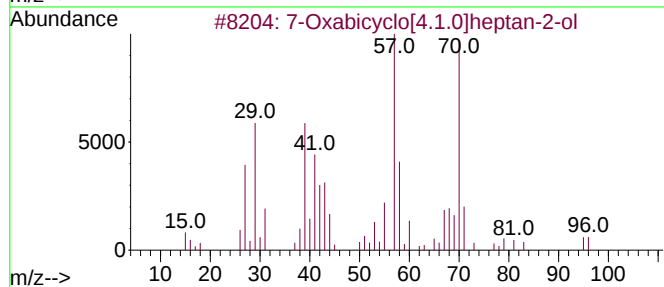
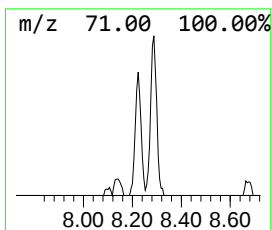
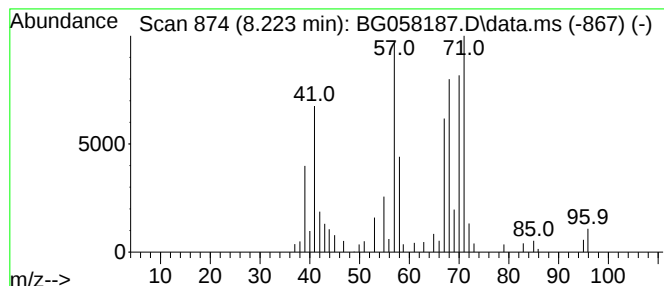
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.223	4.15 ng/ul	56175	1,4-Dichlorobenzene-d4	7.971

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		7-Oxabicyclo[4.1.0]heptan-2-ol	114	C6H10O2	001192-78-5	32
2		Cyclopentanemethanol	100	C6H12O	003637-61-4	27
3		1,4-Pentadiene	68	C5H8	000591-93-5	27
4		1,6-Octadiene, (E)-	110	C8H14	019036-81-8	12
5		2-Butene, 1-methoxy-, (Z)-	86	C5H10O	010034-16-9	10



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Data File : BG058187.D
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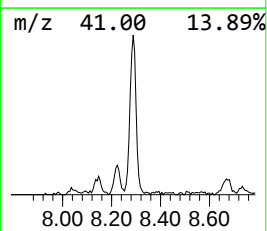
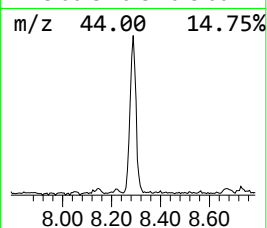
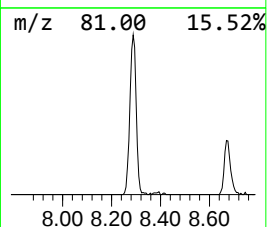
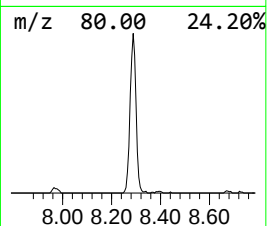
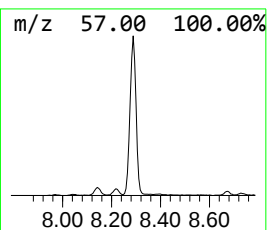
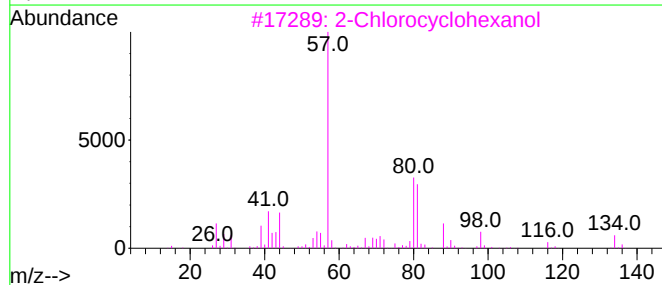
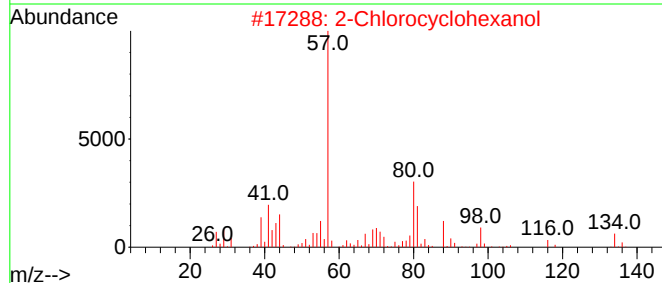
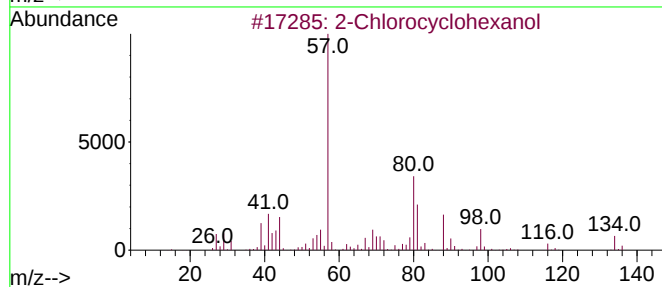
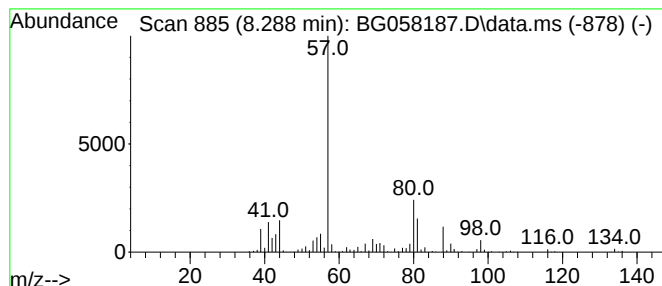
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Chlorocyclohexanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.288	45.29 ng/ul	613661	1,4-Dichlorobenzene-d4	7.971

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	64
4	2-Chlorocyclohexanol			134	C6H11ClO	001561-86-0	64
5	2-Chlorocyclohexanol			134	C6H11ClO	001561-86-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
Data File : BG058187.D
Acq On : 12 Jul 2023 15:11
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Misc :
ALS Vial : 43 Sample Multiplier: 1

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ClientSampleId :
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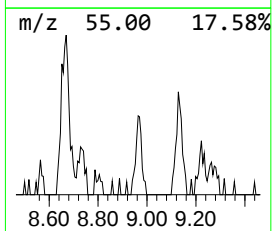
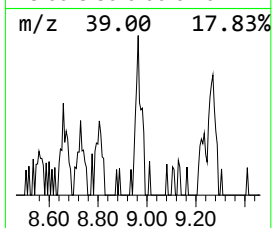
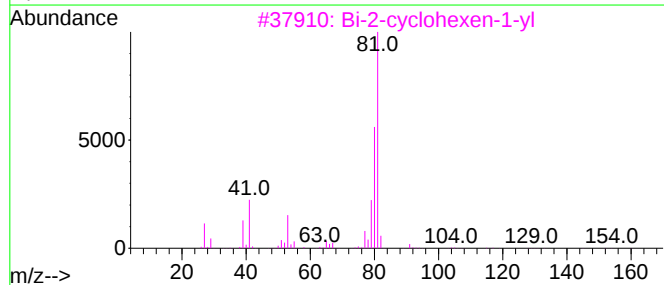
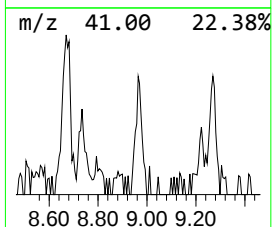
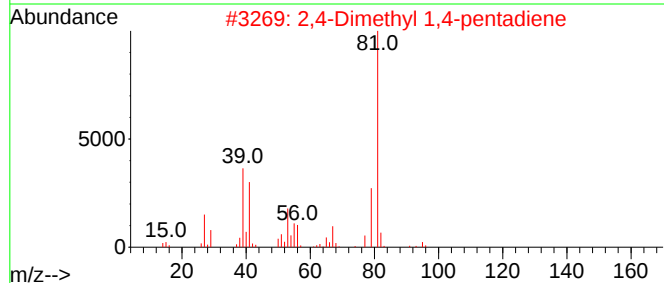
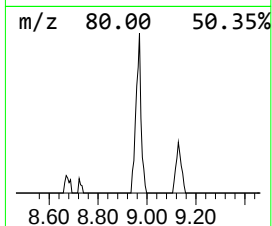
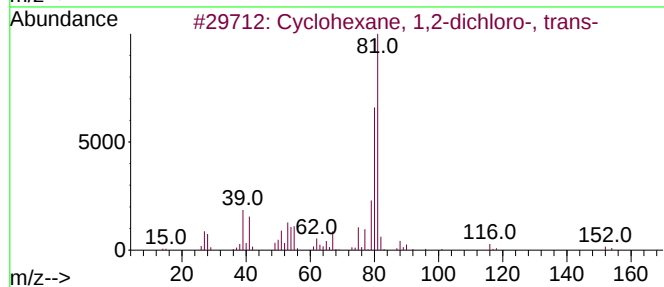
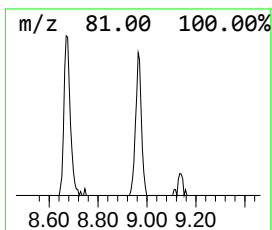
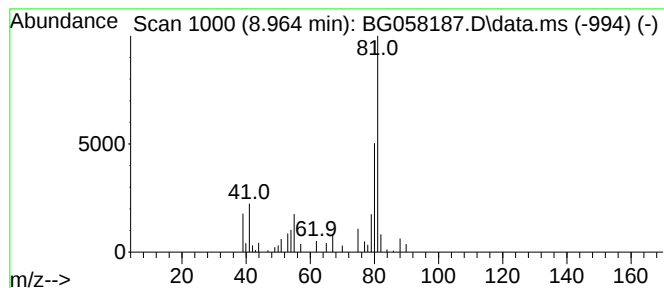
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.964	2.39 ng/ul	32355	1,4-Dichlorobenzene-d4	7.971

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dichloro-, trans-	152	C6H10Cl2	000822-86-6	38
2			2,4-Dimethyl 1,4-pentadiene	96	C7H12	004161-65-3	23
3			Bi-2-cyclohexen-1-yl	162	C12H18	001541-20-4	12
4			1-Penten-3-yne, 2-methyl-	80	C6H8	000926-55-6	9
5			Cyclohexene 3-(tert-butyl)peroxide	170	C10H18O2	051437-25-3	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
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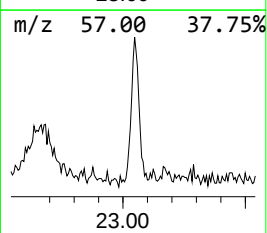
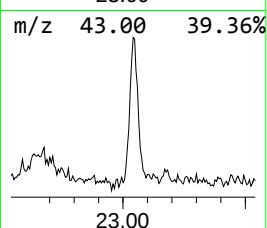
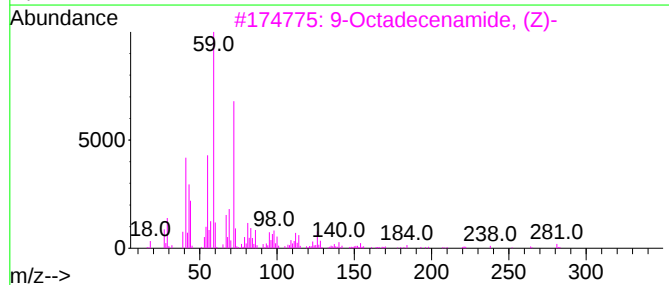
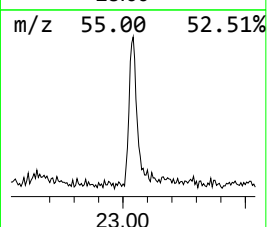
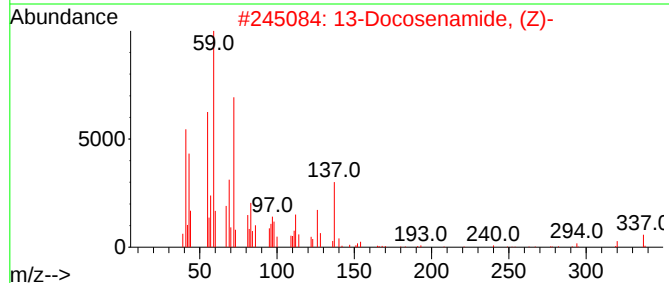
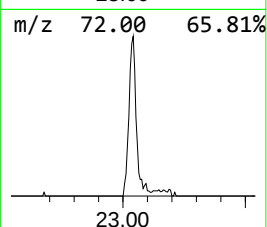
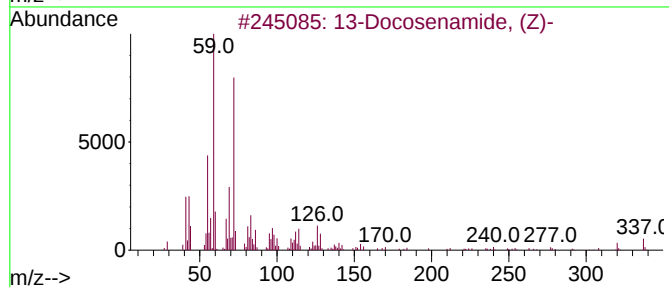
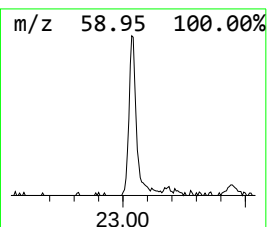
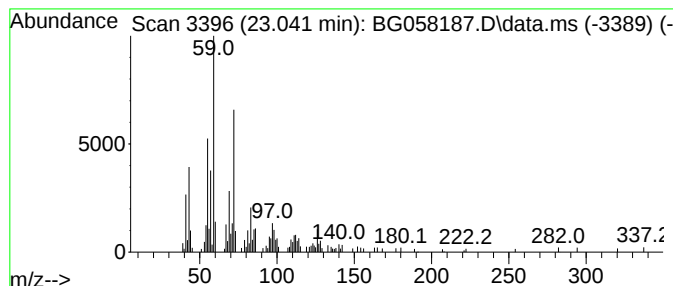
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 13-Docosenamide, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.041	4.04 ng/ul	173893	Chrysene-d12	21.596

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	95
2		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	90
3		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	74
5		Tetradecanamide	227	C14H29NO	000638-58-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
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Acq On : 12 Jul 2023 15:11
Operator : CG/JU
Sample : 03387-11
Misc :
ALS Vial : 43 Sample Multiplier: 1

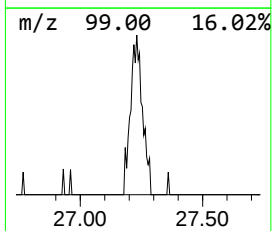
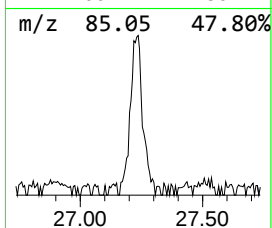
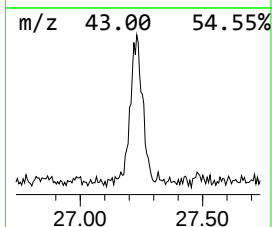
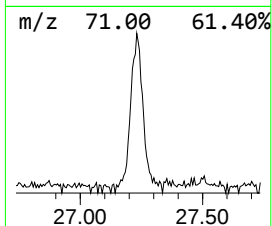
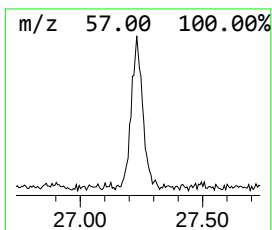
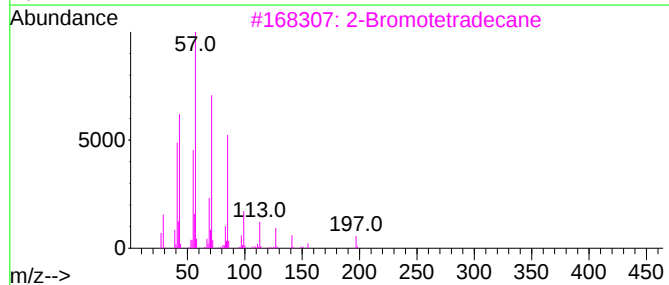
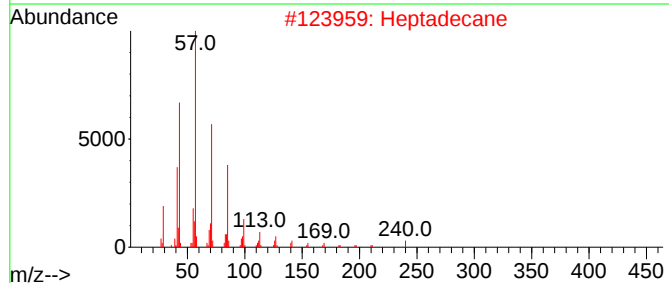
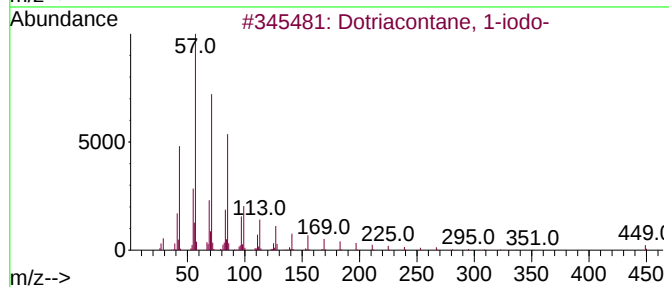
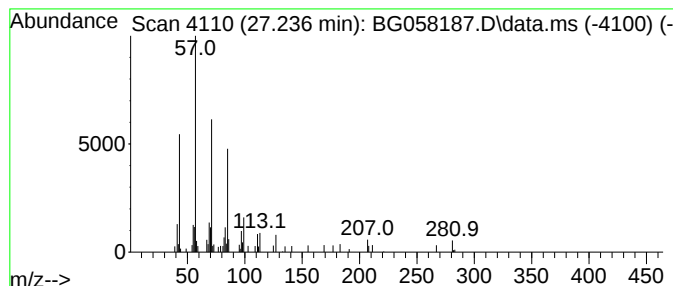
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BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 Dotriacontane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
27.236	2.22 ng/ul	111705	Perylene-d12	24.792	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	80
2	Heptadecane	240	C17H36	000629-78-7	80
3	2-Bromotetradecane	276	C14H29Br	074036-95-6	80
4	Octacosane	394	C28H58	000630-02-4	62
5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
 Data File : BG058187.D
 Acq On : 12 Jul 2023 15:11
 Operator : CG/JU
 Sample : 03387-11
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EX7Z5

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
Butane, 2-metho...	3.141	93.8	ng/ul	1271610	1	7.971	271018	20.0
Cyclohexene	3.217	240.1	ng/ul	3253860	1	7.971	271018	20.0
2-Cyclohexen-1-ol	5.867	8.8	ng/ul	119800	1	7.971	271018	20.0
Cyclohexene, 3-...	6.249	2.6	ng/ul	35602	1	7.971	271018	20.0
2-Cyclohexen-1-one	6.590	18.9	ng/ul	256375	1	7.971	271018	20.0
unknown-01	8.141	3.6	ng/ul	49518	1	7.971	271018	20.0
unknown-02	8.223	4.2	ng/ul	56175	1	7.971	271018	20.0
2-Chlorocyclohe...	8.288	45.3	ng/ul	613661	1	7.971	271018	20.0
unknown-03	8.964	2.4	ng/ul	32355	1	7.971	271018	20.0
13-Docosenamide...	23.041	4.0	ng/ul	173893	5	21.596	861018	20.0
Dotriacontane, ...	27.236	2.2	ng/ul	111705	6	24.792	1006750	20.0