

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV122022\
 Data File : VV029743.D
 Acq On : 20 Dec 2022 12:02
 Operator : SY/MD
 Sample : N6133-01
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EXGJ1

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % 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:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM120722WMA.M
 Title : VOC Analysis

Signal : TIC: VV029743.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.066	3	8	50	rVB	3722821	5010314	100.00%	24.568%
2	1.465	121	132	156	rBV	110239	182119	3.63%	0.893%
3	1.616	174	179	184	rBV2	2231	3085	0.06%	0.015%
4	1.757	219	223	224	rBV2	1386	1120	0.02%	0.005%
5	1.918	261	273	290	rBV	97901	188290	3.76%	0.923%
6	1.989	291	295	300	rVV4	183779	125661	2.51%	0.616%
7	2.150	340	345	352	rBV5	630441	532717	10.63%	2.612%
8	2.950	586	594	598	rBV	24276	34760	0.69%	0.170%
9	2.989	599	606	609	rVV	49191	61614	1.23%	0.302%
10	3.031	609	619	637	rVB9	1120957	2055693	41.03%	10.080%
11	3.445	740	748	753	rBV7	1045689	1596420	31.86%	7.828%
12	3.796	855	857	858	rBV2	3488	1410	0.03%	0.007%
13	5.310	1317	1328	1342	rBV	148474	283984	5.67%	1.393%
14	5.802	1470	1481	1494	rBV2	304200	590442	11.78%	2.895%
15	6.211	1601	1608	1618	rBV	40338	54772	1.09%	0.269%
16	6.323	1619	1643	1674	rVB	322738	898023	17.92%	4.403%
17	6.551	1705	1714	1728	rVB2	38537	62635	1.25%	0.307%
18	6.690	1742	1757	1776	rBV	308339	576795	11.51%	2.828%
19	6.815	1785	1796	1826	rVB	26155	67772	1.35%	0.332%
20	7.088	1875	1881	1884	rBV4	1169	1298	0.03%	0.006%
21	7.191	1901	1913	1917	rBV4	4188	7276	0.15%	0.036%
22	7.211	1917	1919	1923	rVB4	1184	958	0.02%	0.005%
23	7.403	1964	1979	2005	rBV	153123	274353	5.48%	1.345%
24	7.603	2032	2041	2051	rBV9	3243	6666	0.13%	0.033%
25	7.683	2052	2066	2081	rBV	538350	921631	18.39%	4.519%
26	7.928	2130	2142	2167	rBV2	108164	192889	3.85%	0.946%
27	8.104	2187	2197	2199	rBV7	997	1239	0.02%	0.006%
28	8.169	2199	2217	2236	rBV2	49784	93803	1.87%	0.460%
29	8.333	2248	2268	2302	rBV	282593	750082	14.97%	3.678%
30	8.468	2306	2310	2328	rVB8	8539	16904	0.34%	0.083%
31	8.670	2370	2373	2379	rVB6	1063	969	0.02%	0.005%
32	8.847	2421	2428	2437	rVB9	1663	2656	0.05%	0.013%
33	8.911	2446	2448	2454	rVB4	790	745	0.01%	0.004%
34	9.030	2468	2485	2521	rBV	603378	1074945	21.45%	5.271%
35	9.178	2526	2531	2534	rVB7	951	925	0.02%	0.005%
36	9.198	2534	2537	2545	rVB8	1043	1416	0.03%	0.007%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM120722WMA.M
 Title : VOC Analysis

37	9.288	2557	2565	2571	rBV6	1842	2617	0.05%	0.013%
38	9.378	2589	2593	2597	rVB5	1268	1160	0.02%	0.006%
39	9.455	2611	2617	2621	rBV5	1195	1526	0.03%	0.007%
40	9.574	2648	2654	2656	rBV6	1739	1673	0.03%	0.008%
41	9.699	2691	2693	2702	rVB6	1478	1660	0.03%	0.008%
42	9.802	2720	2725	2731	rBV4	2666	2915	0.06%	0.014%
43	9.899	2748	2755	2776	rVB10	5536	10722	0.21%	0.053%
44	10.034	2790	2797	2802	rVB7	901	1273	0.03%	0.006%
45	10.075	2806	2810	2814	rBV5	1110	839	0.02%	0.004%
46	10.284	2853	2875	2891	rBV	398060	650476	12.98%	3.190%
47	10.419	2906	2917	2933	rBV	38823	65444	1.31%	0.321%
48	10.516	2943	2947	2954	rVB6	905	1052	0.02%	0.005%
49	10.564	2954	2962	2967	rBV6	1971	2892	0.06%	0.014%
50	10.587	2967	2969	2975	rVB4	943	882	0.02%	0.004%
51	10.635	2978	2984	2993	rVB8	2395	3525	0.07%	0.017%
52	10.802	3032	3036	3043	rBV7	1737	2417	0.05%	0.012%
53	10.853	3045	3052	3055	rBV7	2609	3160	0.06%	0.015%
54	10.969	3070	3088	3103	rVB9	8565	21962	0.44%	0.108%
55	11.027	3103	3106	3109	rBV5	994	840	0.02%	0.004%
56	11.075	3112	3121	3129	rBV7	3127	5643	0.11%	0.028%
57	11.178	3144	3153	3162	rBV7	5332	8921	0.18%	0.044%
58	11.217	3162	3165	3173	rVB6	1360	1793	0.04%	0.009%
59	11.284	3173	3186	3207	rBV	673991	1111688	22.19%	5.451%
60	11.548	3251	3268	3288	rBV	304672	552253	11.02%	2.708%
61	11.651	3289	3300	3320	rVB	649004	1040597	20.77%	5.103%
62	11.992	3395	3406	3408	rBV4	8991	12441	0.25%	0.061%
63	12.156	3455	3457	3463	rVB6	1007	884	0.02%	0.004%
64	12.249	3472	3486	3499	rBV3	17847	33408	0.67%	0.164%
65	12.349	3509	3517	3525	rBV9	5509	9345	0.19%	0.046%
66	12.439	3540	3545	3549	rBV7	1901	2210	0.04%	0.011%
67	12.651	3609	3611	3621	rVB9	2049	1959	0.04%	0.010%
68	12.725	3629	3634	3643	rVB7	2428	3303	0.07%	0.016%
69	12.792	3647	3655	3656	rBV6	5488	5919	0.12%	0.029%
70	12.844	3665	3671	3682	rVB6	14375	26019	0.52%	0.128%
71	12.956	3698	3706	3715	rBV6	7671	10570	0.21%	0.052%
72	13.101	3743	3751	3759	rBV6	10042	17015	0.34%	0.083%
73	13.233	3783	3792	3795	rBV9	1715	2780	0.06%	0.014%
74	13.284	3796	3808	3812	rBV10	6211	12769	0.25%	0.063%
75	13.358	3827	3831	3837	rVB9	2721	2788	0.06%	0.014%
76	13.455	3847	3861	3865	rBV9	9919	20638	0.41%	0.101%

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM120722WMA.M
 Title : VOC Analysis

77	13.500	3868	3875	3887	rVB	18791	35935	0.72%	0.176%
78	13.622	3910	3913	3921	rVB9	4488	5443	0.11%	0.027%
79	13.953	4013	4016	4017	rBV3	1136	680	0.01%	0.003%
80	14.027	4032	4039	4048	rBV2	7126	12110	0.24%	0.059%
81	14.107	4058	4064	4073	rBV2	4673	7821	0.16%	0.038%
82	14.320	4121	4130	4138	rBV6	6562	10180	0.20%	0.050%
83	14.426	4152	4163	4169	rBV6	5691	12308	0.25%	0.060%
84	14.651	4231	4233	4236	rBV4	1455	1177	0.02%	0.006%
85	14.805	4276	4281	4288	rVB9	5682	6888	0.14%	0.034%
86	14.892	4302	4308	4317	rVB10	3262	5230	0.10%	0.026%
87	14.940	4320	4323	4324	rBV3	1538	894	0.02%	0.004%
88	14.988	4335	4338	4343	rBV5	1369	1407	0.03%	0.007%
89	15.024	4343	4349	4350	rBV6	2687	2361	0.05%	0.012%
90	15.130	4380	4382	4386	rVB5	1631	866	0.02%	0.004%
91	15.162	4387	4392	4393	rBV5	1119	864	0.02%	0.004%
92	15.197	4401	4403	4406	rVB4	1474	734	0.01%	0.004%
93	15.255	4413	4421	4422	rBV6	4523	4987	0.10%	0.024%
94	15.368	4446	4456	4462	rBV8	13544	23937	0.48%	0.117%
95	15.474	4487	4489	4494	rBV6	1381	1373	0.03%	0.007%
96	15.760	4570	4578	4589	rVB2	14533	20684	0.41%	0.101%
97	15.930	4620	4631	4661	rBV	390246	706723	14.11%	3.465%
98	17.082	4981	4989	4999	rBV2	29900	43971	0.88%	0.216%
99	17.429	5091	5097	5115	rVB4	14500	31182	0.62%	0.153%
100	17.602	5140	5151	5165	rBV2	59263	119339	2.38%	0.585%

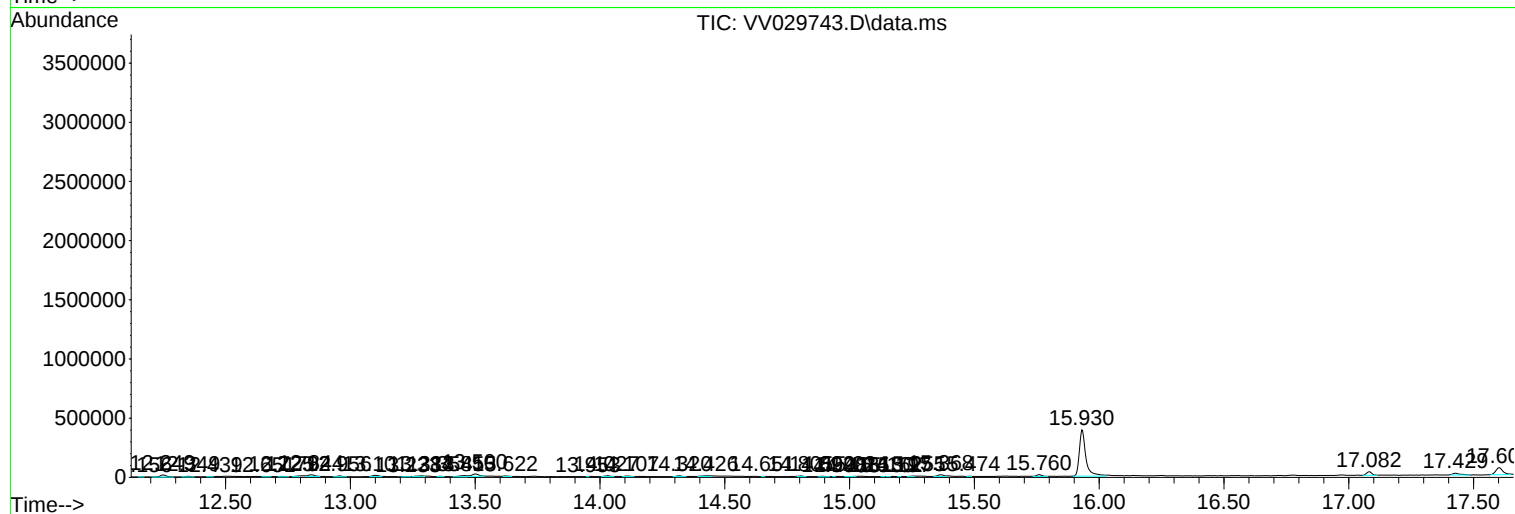
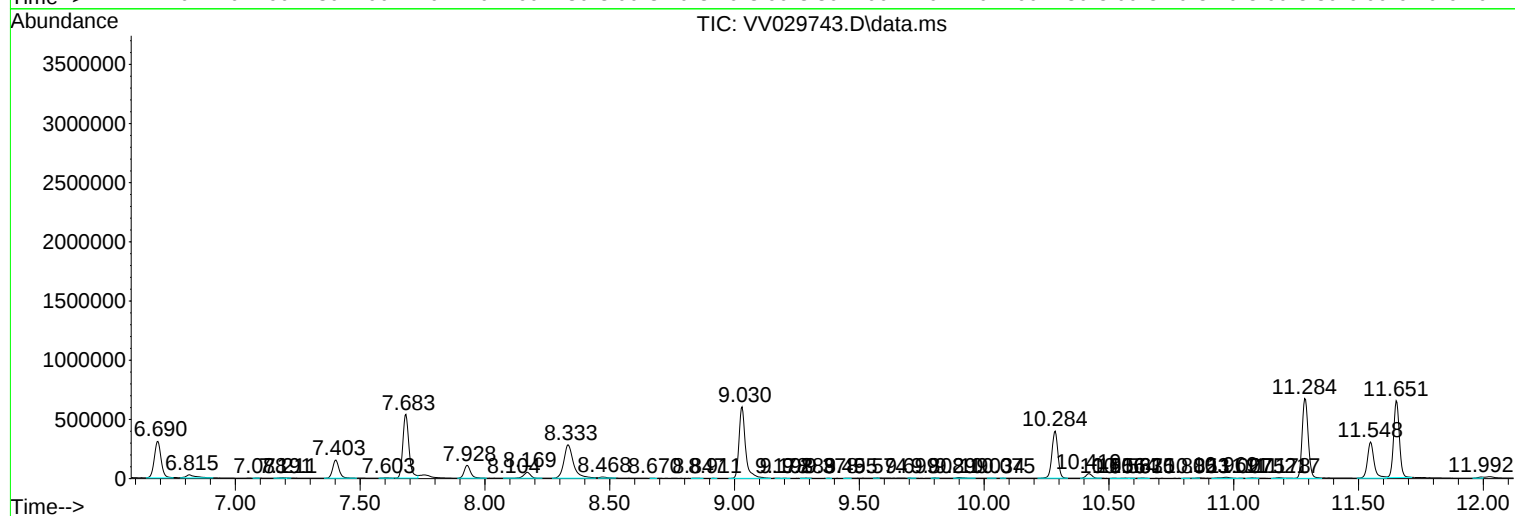
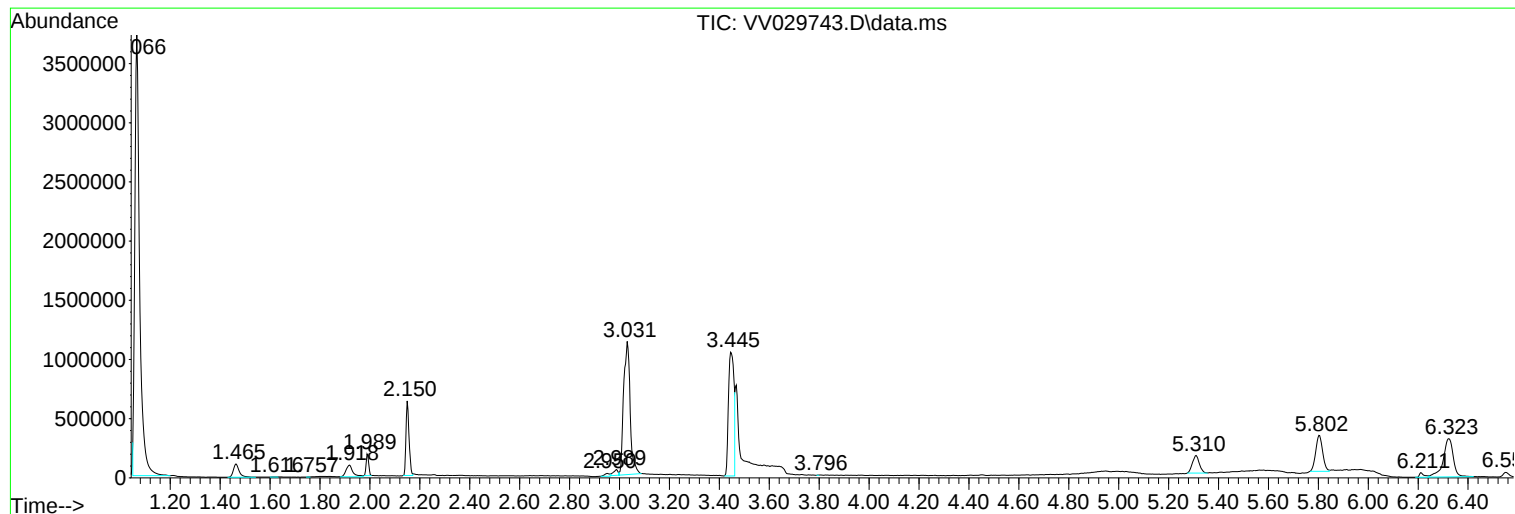
Sum of corrected areas: 20393453

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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXGJ1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM120722WMA.M
Quant Title : VOC Analysis

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



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Data File : VV029743.D
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Instrument :
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ClientSampleId :
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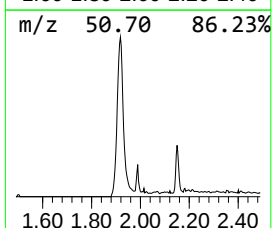
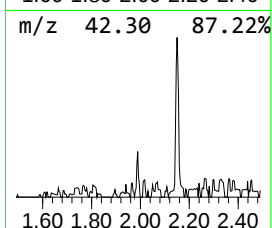
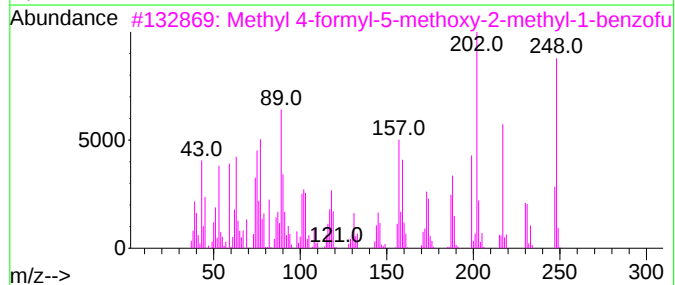
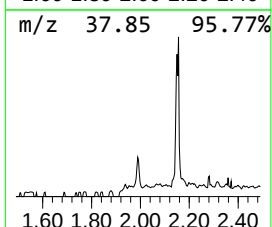
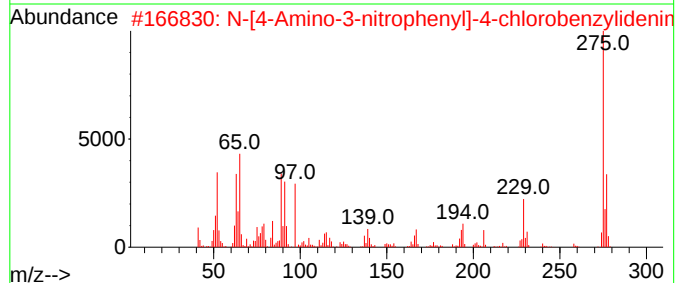
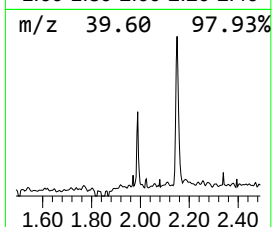
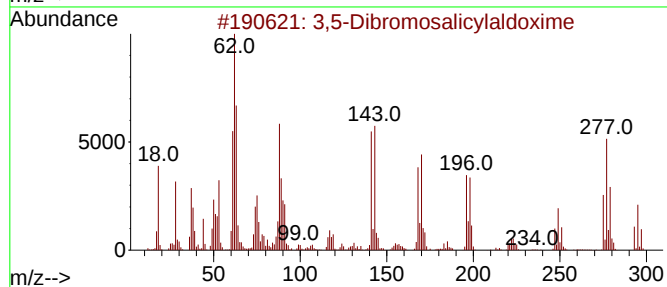
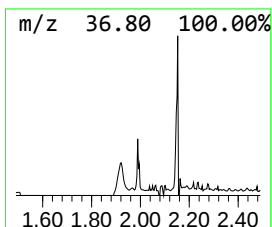
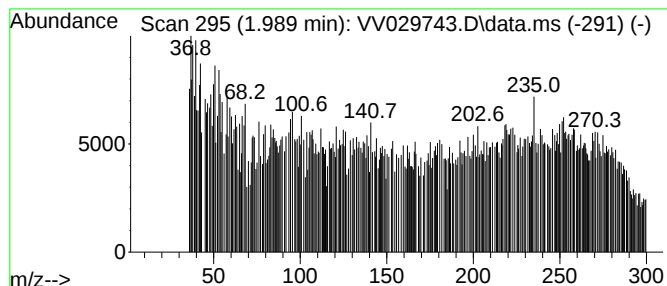
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM120722WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 3,5-Dibromosalicylaldehyde Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.989	7.00 ug/L	125661	1,4-Difluorobenzene	6.323

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Dibromosalicylaldehyde	293	C7H5Br2NO2	021386-43-6	90
2			N-[4-Amino-3-nitrophenyl]-4-chlo...	275	C13H10ClN3O2	1000213-83-4	55
3			Methyl 4-formyl-5-methoxy-2-meth...	248	C13H12O5	1000410-12-5	55
4			3-Bromo-5-phenyl-1H-1,2,4-triazole	223	C8H6BrN3	015777-59-0	45
5			Methylarsine dibromide	248	CH3AsBr2	000676-70-0	38



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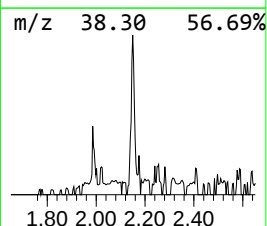
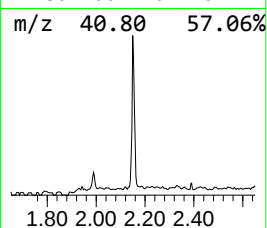
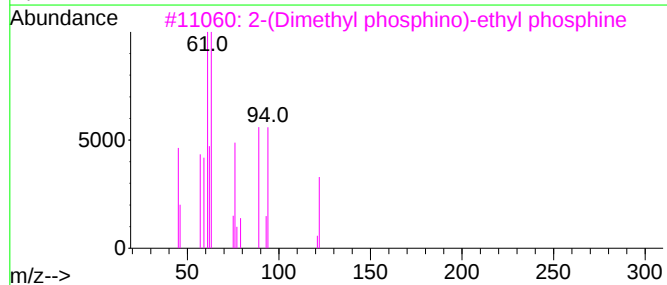
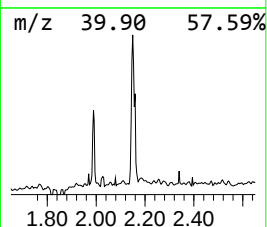
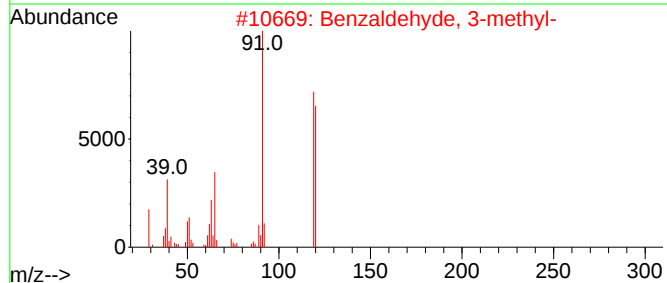
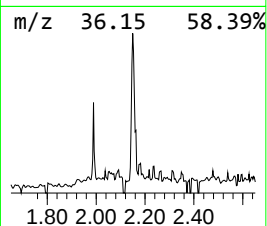
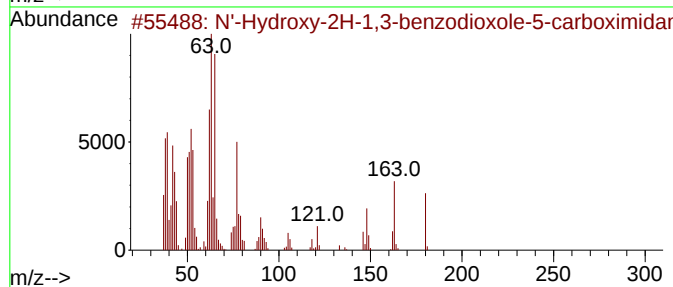
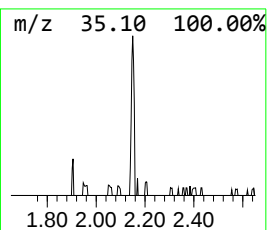
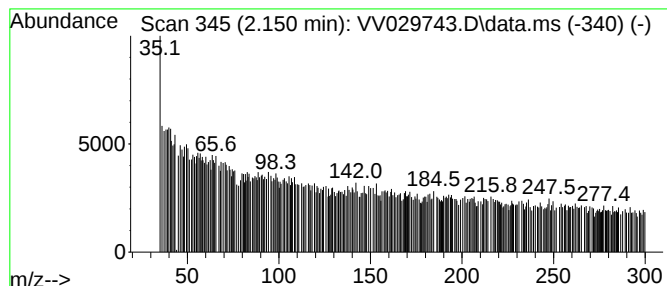
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM120722WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 N'-Hydroxy-2H-1,3-benzodiox... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.150	29.66 ug/L	532717	1,4-Difluorobenzene	6.323

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N'-Hydroxy-2H-1,3-benzodioxole-5...	180	C8H8N2O3	004720-72-3	95
2			Benzaldehyde, 3-methyl-	120	C8H8O	000620-23-5	50
3			2-(Dimethyl phosphino)-ethyl pho...	122	C4H12P2	054772-64-4	47
4			2,4,6-Cycloheptatrien-1-one, 3-b...	200	C7H5BrO2	004584-68-3	46
5			3-pyridinesulfonyl chloride, 5,6...	245	C5H2Cl3N02S	1000401-30-4	44



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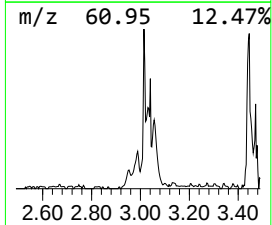
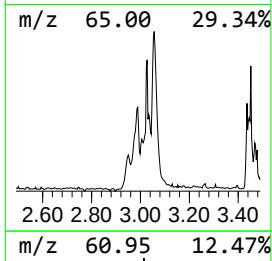
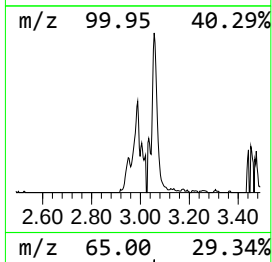
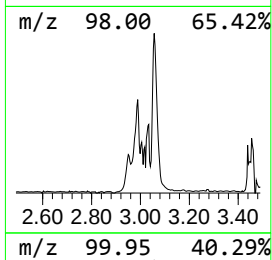
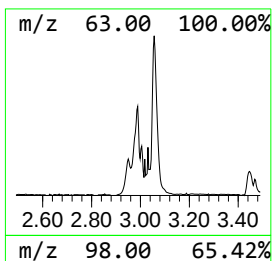
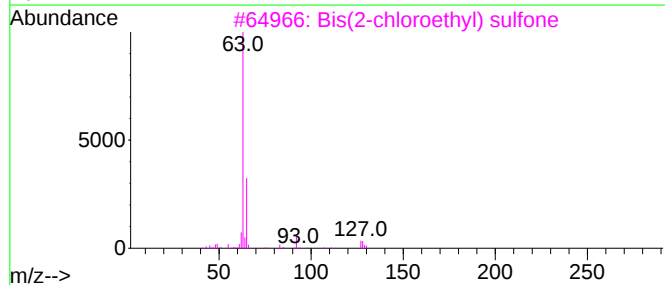
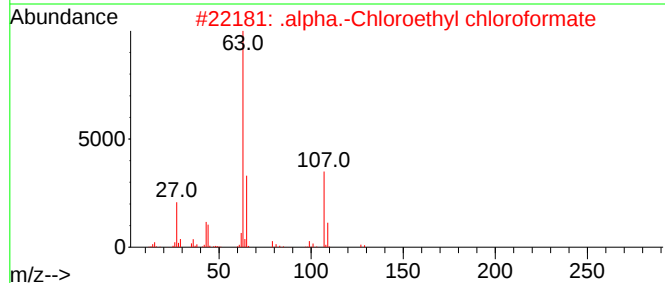
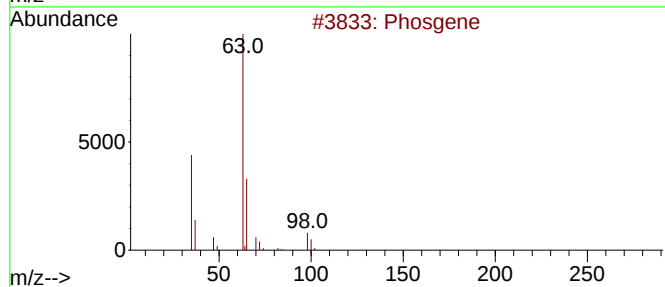
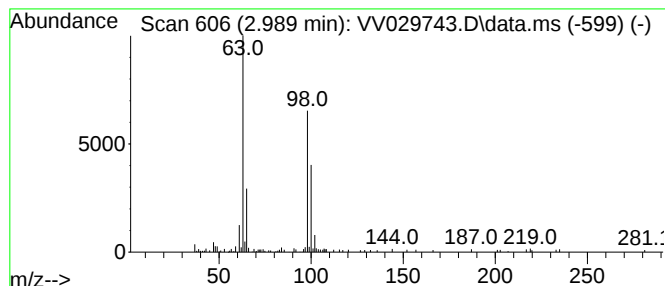
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Quant Title : VOC Analysis

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

Peak Number 4 Phosgene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.989	3.43 ug/L	61614	1,4-Difluorobenzene	6.323

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phosgene	98	CCl2O	000075-44-5	64
2			.alpha.-Chloroethyl chloroformate	142	C3H4Cl2O2	050893-53-3	45
3			Bis(2-chloroethyl) sulfone	190	C4H8Cl2O2S	000471-03-4	45
4			Ethene, (methylsulfonyl)-	106	C3H6O2S	003680-02-2	39
5			Ethane, 1,1-dichloro-	98	C2H4Cl2	000075-34-3	39



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV122022\
Data File : VV029743.D
Acq On : 20 Dec 2022 12:02
Operator : SY/MD
Sample : N6133-01
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXGJ1

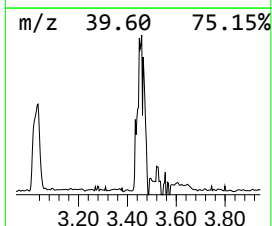
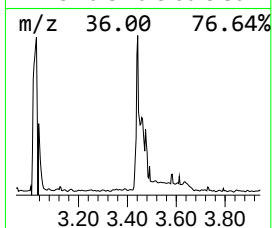
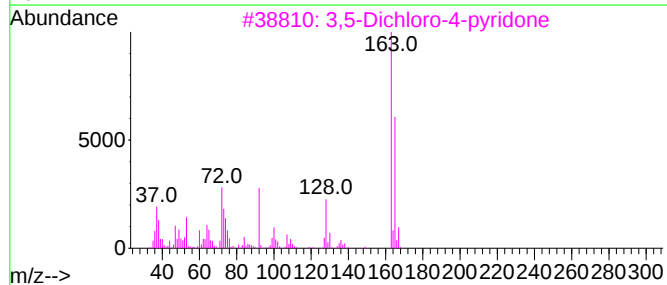
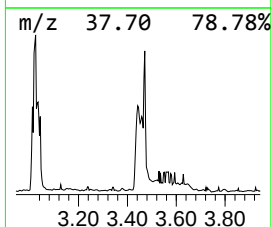
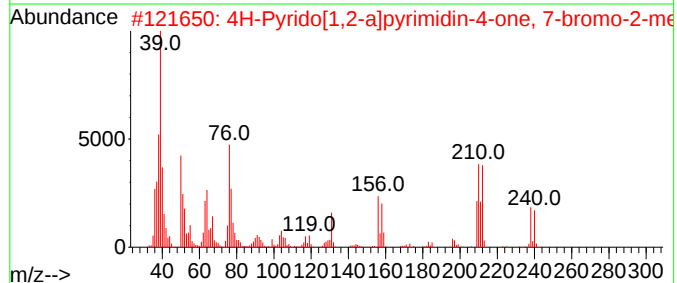
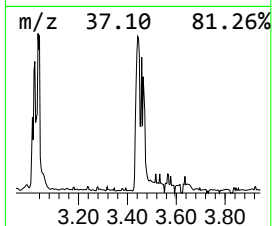
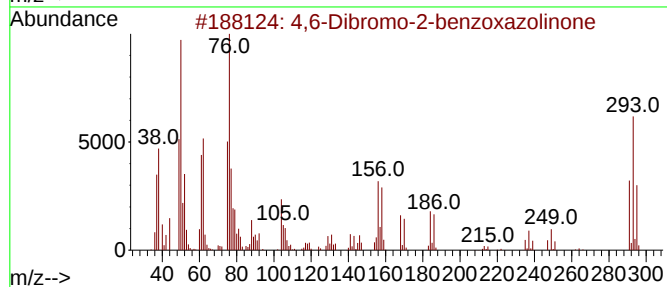
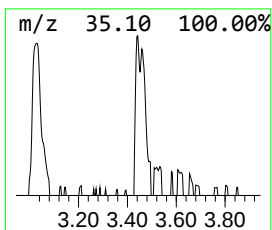
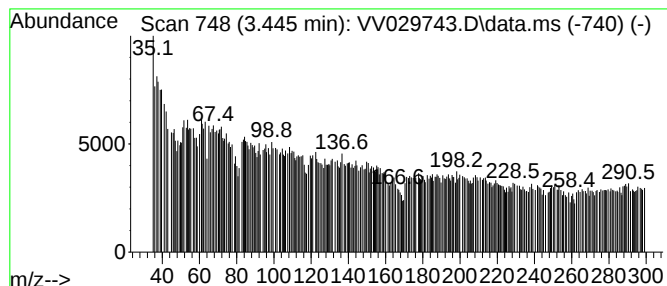
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Quant Title : VOC Analysis

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

Peak Number 5 4,6-Dibromo-2-benzoxazolinone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.445	88.89 ug/L	1596420	1,4-Difluorobenzene	6.323

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,6-Dibromo-2-benzoxazolinone	291	C7H3Br2NO2	1000260-92-6	74
2			4H-Pyrido[1,2-a]pyrimidin-4-one,...	238	C9H7BrN2O	064500-11-4	59
3			3,5-Dichloro-4-pyridone	163	C5H3Cl2NO	017228-70-5	55
4			5-Chloro-3-phenyl-1,2,4-oxadiazole	180	C8H5ClN2O	000827-44-1	55
5			1,2,5-Oxadiazole-3-carbohydrazon...	275	C10H9N7O3	1000276-22-7	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV122022\
Data File : VV029743.D
Acq On : 20 Dec 2022 12:02
Operator : SY/MD
Sample : N6133-01
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXGJ1

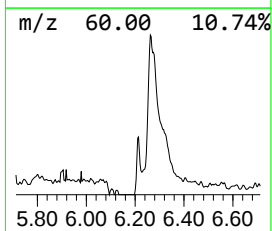
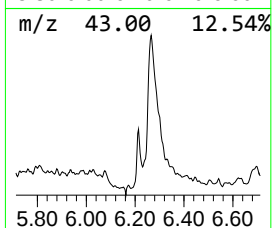
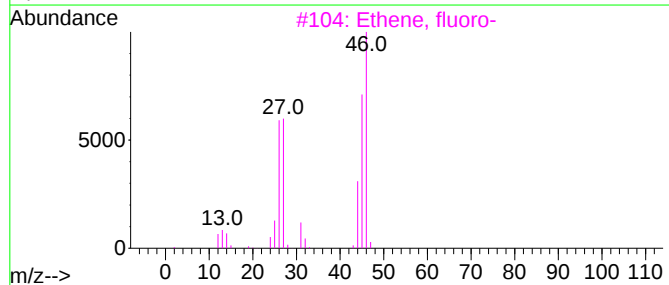
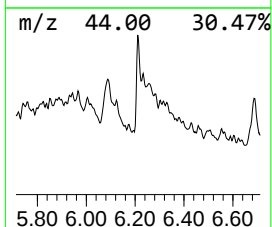
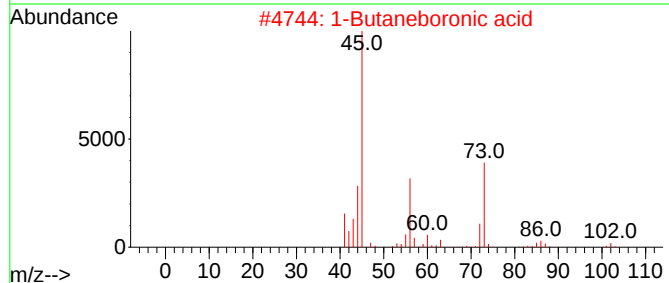
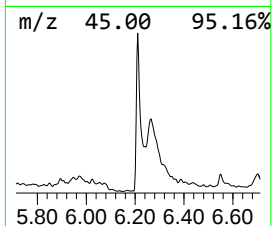
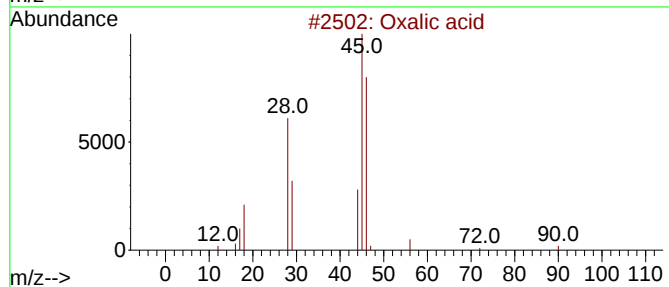
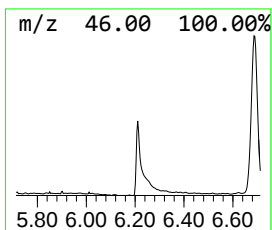
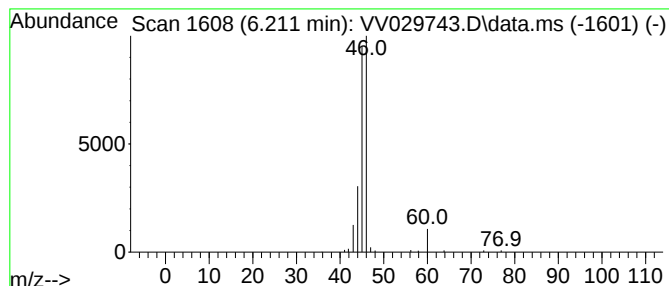
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Quant Title : VOC Analysis

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

Peak Number 6 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.211	3.05 ug/L	54772	1,4-Difluorobenzene	6.323

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxalic acid	90	C2H2O4	000144-62-7	43
2		1-Butaneboronic acid	102	C4H11BO2	004426-47-5	25
3		Ethene, fluoro-	46	C2H3F	000075-02-5	9
4		Formic acid	46	CH2O2	000064-18-6	9
5		Formic acid	46	CH2O2	000064-18-6	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV122022\
Data File : VV029743.D
Acq On : 20 Dec 2022 12:02
Operator : SY/MD
Sample : N6133-01
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXGJ1

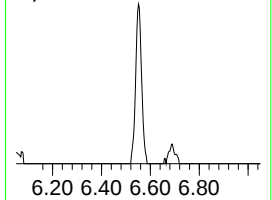
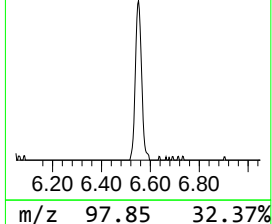
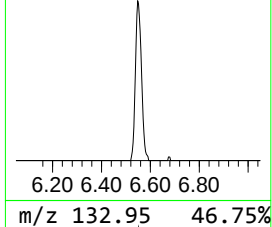
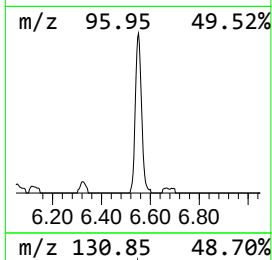
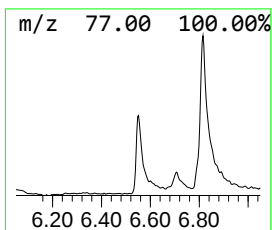
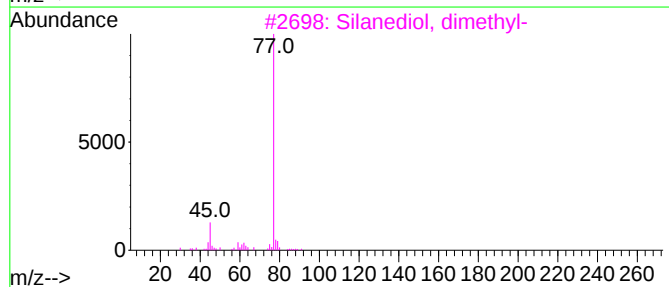
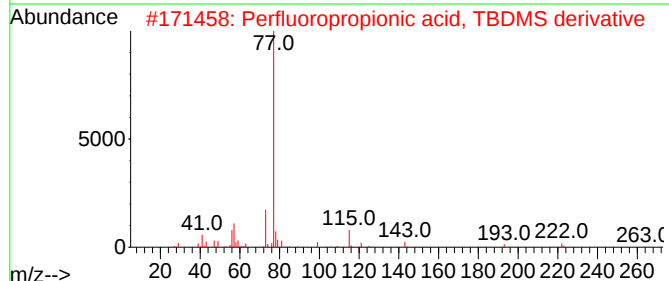
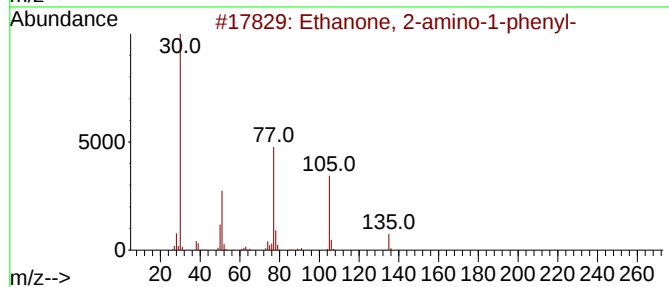
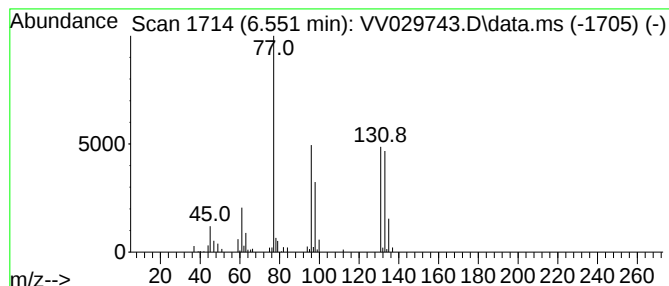
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Quant Title : VOC Analysis

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

Peak Number 7 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.551	3.49 ug/L	62635	1,4-Difluorobenzene	6.323

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 2-amino-1-phenyl-	135	C8H9NO	000613-89-8	12
2			Perfluoropropionic acid, TBDMS d...	278	C9H15F5O2Si	1010376-29-3	10
3			Silanediol, dimethyl-	92	C2H8O2Si	001066-42-8	9
4			Acetamide, 2-phenoxy-N-(4-benzoy...	346	C21H18N2O3	312755-44-5	9
5			Silanediol, dimethyl-	92	C2H8O2Si	001066-42-8	8



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV122022\
Data File : VV029743.D
Acq On : 20 Dec 2022 12:02
Operator : SY/MD
Sample : N6133-01
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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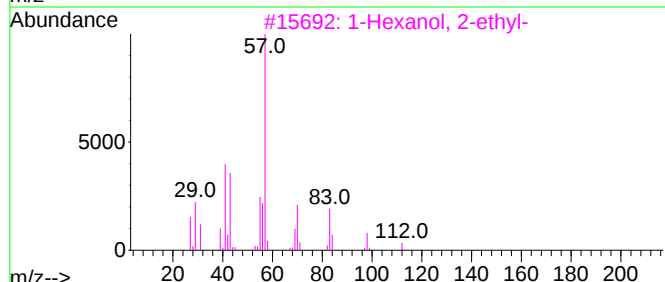
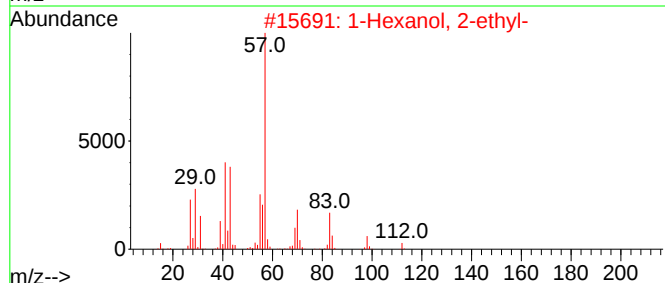
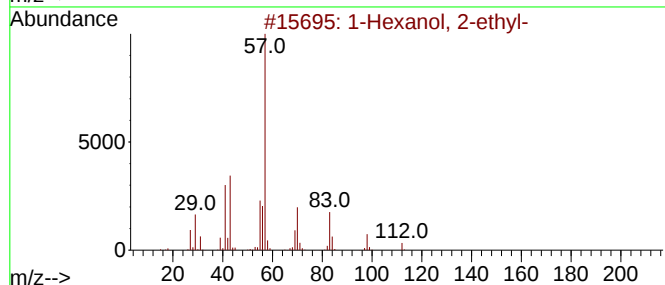
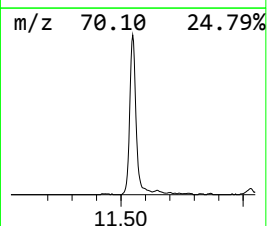
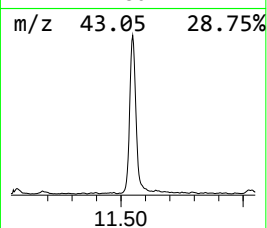
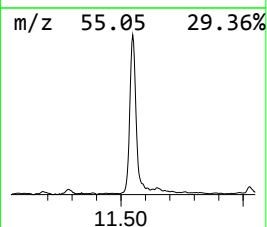
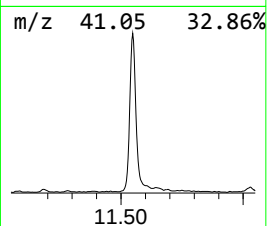
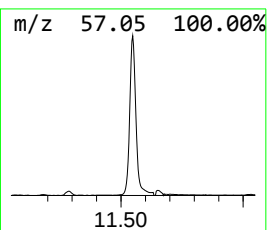
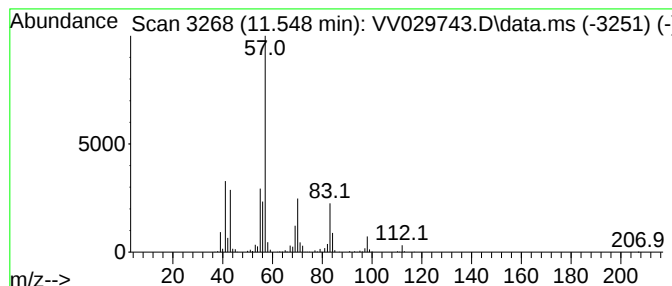
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Quant Title : VOC Analysis

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

Peak Number 12 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.548	24.84 ug/L	552253	1,4-Dichlorobenzene-d4	11.284

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	90
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	90
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	90
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV122022\
Data File : VV029743.D
Acq On : 20 Dec 2022 12:02
Operator : SY/MD
Sample : N6133-01
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXGJ1

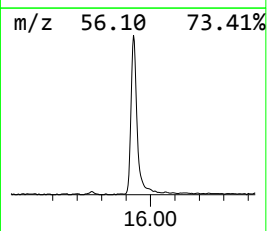
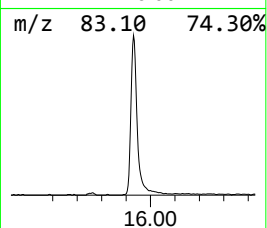
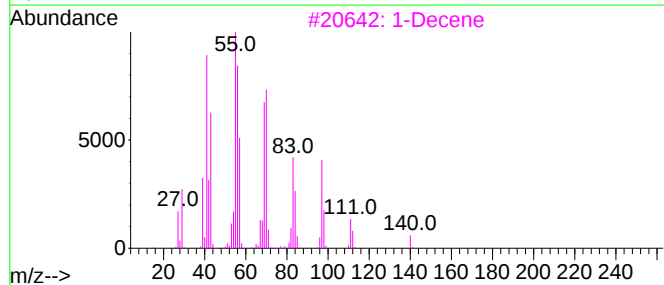
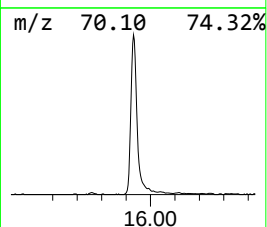
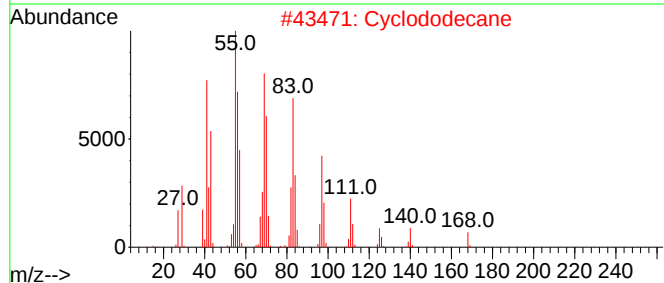
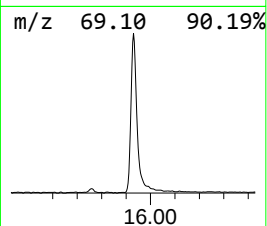
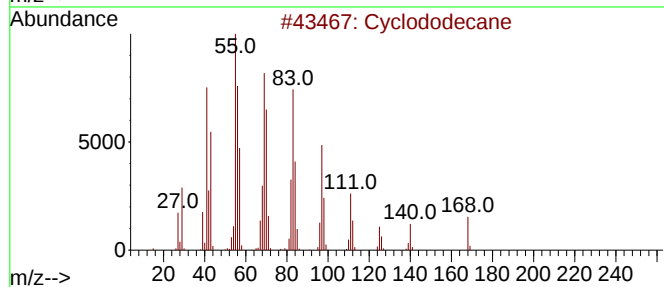
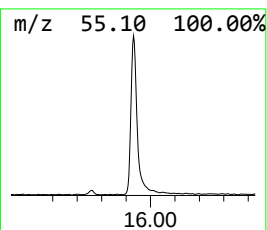
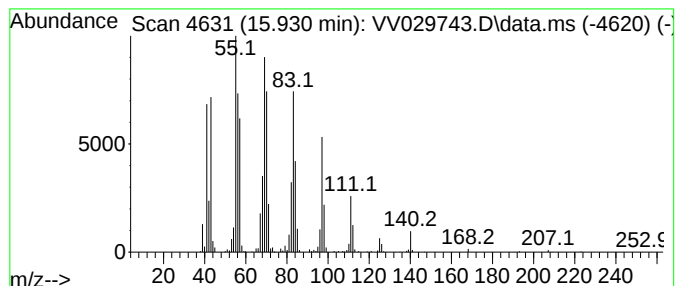
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Quant Title : VOC Analysis

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

Peak Number 13 (DEL) Alkane: Cyclic15.930 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.930	31.79 ug/L	706723	1,4-Dichlorobenzene-d4	11.284

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclododecane	168	C12H24	000294-62-2	97
2		Cyclododecane	168	C12H24	000294-62-2	96
3		1-Decene	140	C10H20	000872-05-9	95
4		Cyclopropane, nonyl-	168	C12H24	074663-85-7	95
5		1-Dodecanol	186	C12H26O	000112-53-8	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV122022\
Data File : VV029743.D
Acq On : 20 Dec 2022 12:02
Operator : SY/MD
Sample : N6133-01
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXGJ1

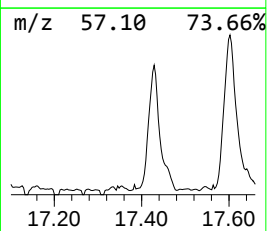
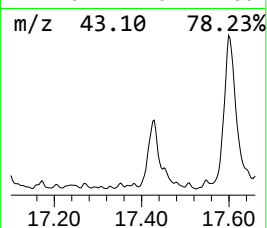
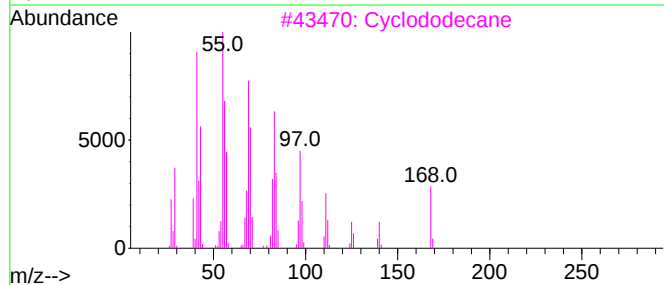
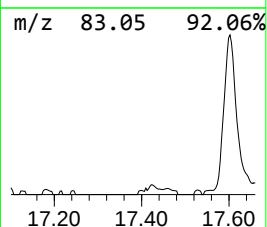
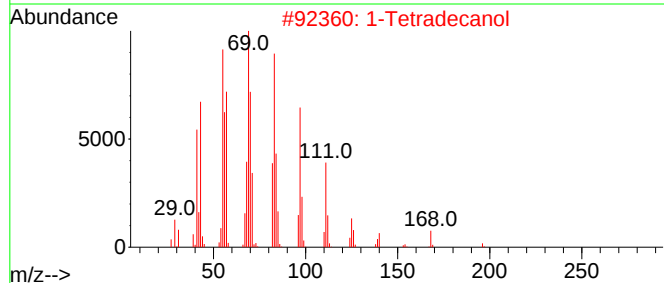
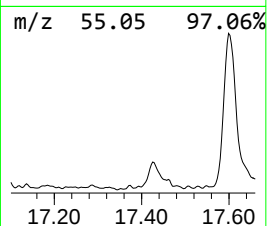
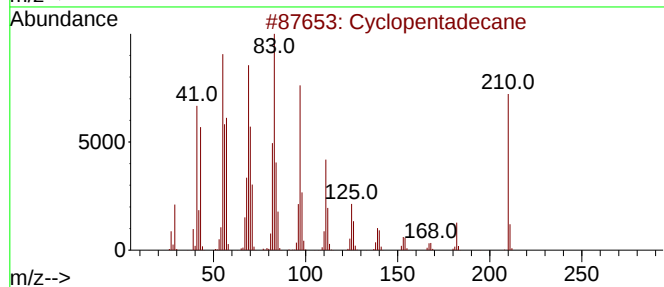
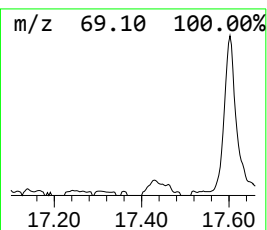
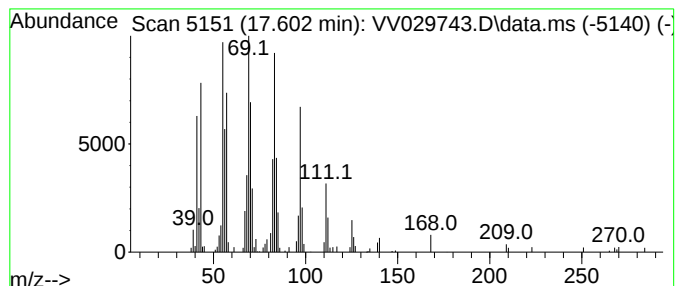
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Quant Title : VOC Analysis

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

Peak Number 14 (DEL) Alkane: Cyclic17.602 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.602	5.37 ug/L	119339	1,4-Dichlorobenzene-d4	11.284

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	98
2			1-Tetradecanol	214	C14H30O	000112-72-1	95
3			Cyclododecane	168	C12H24	000294-62-2	94
4			Cyclododecane	168	C12H24	000294-62-2	93
5			n-Pentadecanol	228	C15H32O	000629-76-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW122022\
Data File : VW029743.D
Acq On : 20 Dec 2022 12:02
Operator : SY/MD
Sample : N6133-01
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
EXGJ1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM120722WMA.M
Quant Title : VOC Analysis

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3,5-Dibromosali...	1.989	7.0	ug/L	125661	1	6.323	898023	50.0
N'-Hydroxy-2H-1...	2.150	29.7	ug/L	532717	1	6.323	898023	50.0
Phosgene	2.989	3.4	ug/L	61614	1	6.323	898023	50.0
4,6-Dibromo-2-b...	3.445	88.9	ug/L	1596420	1	6.323	898023	50.0
unknown-01	6.211	3.0	ug/L	54772	1	6.323	898023	50.0
unknown-02	6.551	3.5	ug/L	62635	1	6.323	898023	50.0
1-Hexanol, 2-et...	11.548	24.8	ug/L	552253	3	11.284	1111690	50.0
(DEL) Alkane: C...	15.930	31.8	ug/L	706723	3	11.284	1111690	50.0
(DEL) Alkane: C...	17.602	5.4	ug/L	119339	3	11.284	1111690	50.0