

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101921\
 Data File : VW020628.D
 Acq On : 20 Oct 2021 00:25
 Operator : SY/VA
 Sample : M4263-11
 Misc : 6.10g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 GB7F6

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_W\Method\SFAMWLM100821SMA.M
 Title : SFAM01.0

Signal : TIC: VW020628.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.343	67	72	83	rVB	72816	124989	13.48%	1.683%
2	2.489	91	96	106	rVB	22393	39463	4.25%	0.531%
3	2.806	145	148	152	rVB	1204	2057	0.22%	0.028%
4	2.879	152	160	168	rBV	26224	56652	6.11%	0.763%
5	3.019	178	183	190	rVB4	2119	4913	0.53%	0.066%
6	3.294	221	228	232	rBV5	897	2037	0.22%	0.027%
7	3.379	233	242	252	rVB	8794	21271	2.29%	0.286%
8	3.531	265	267	269	rBV2	289	373	0.04%	0.005%
9	3.720	291	298	299	rBV2	470	905	0.10%	0.012%
10	3.836	312	317	320	rBV3	384	569	0.06%	0.008%
11	3.903	323	328	333	rVV4	281	594	0.06%	0.008%
12	4.013	333	346	357	rVV2	96979	304906	32.87%	4.105%
13	4.123	358	364	380	rVV	23742	66749	7.20%	0.899%
14	4.245	381	384	395	rVB5	642	1672	0.18%	0.023%
15	4.379	397	406	417	rBV4	1118	3710	0.40%	0.050%
16	4.616	441	445	447	rBV	235	342	0.04%	0.005%
17	4.684	447	456	462	rBV2	1603	4301	0.46%	0.058%
18	4.909	484	493	500	rVV3	3472	11456	1.24%	0.154%
19	5.007	503	509	520	rVV3	4753	15084	1.63%	0.203%
20	5.440	576	580	585	rVV3	384	646	0.07%	0.009%
21	5.787	627	637	652	rVB3	4578	14673	1.58%	0.198%
22	5.939	652	662	670	rVB2	3011	8791	0.95%	0.118%
23	6.122	683	692	701	rVV4	805	2367	0.26%	0.032%
24	6.897	809	819	830	rBV2	15226	42421	4.57%	0.571%
25	7.079	839	849	861	rBV	30065	83618	9.02%	1.126%
26	7.177	861	865	879	rVB	3017	8385	0.90%	0.113%
27	7.646	931	942	960	rBV	154791	377437	40.69%	5.082%
28	8.274	1035	1045	1059	rBV2	309117	927490	100.00%	12.488%
29	8.549	1085	1090	1107	rVB4	10675	31952	3.44%	0.430%
30	8.707	1107	1116	1125	rBV5	5413	14517	1.57%	0.195%
31	8.841	1129	1138	1154	rBV	281103	545230	58.79%	7.341%
32	9.036	1165	1170	1171	rBV3	2026	2876	0.31%	0.039%
33	9.158	1186	1190	1196	rVB3	385	654	0.07%	0.009%
34	9.274	1200	1209	1224	rBV	227196	453179	48.86%	6.102%
35	9.408	1224	1231	1244	rVB	60476	113106	12.19%	1.523%
36	9.506	1244	1247	1251	rBV2	266	413	0.04%	0.006%

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 Title : SFAM01.0

37	9.664	1268	1273	1279	rVB3	502	809	0.09%	0.011%
38	9.920	1311	1315	1319	rBV2	725	1201	0.13%	0.016%
39	10.036	1325	1334	1346	rBV	114176	205030	22.11%	2.761%
40	10.128	1346	1349	1354	rVB3	490	671	0.07%	0.009%
41	10.213	1357	1363	1370	rBV3	476	925	0.10%	0.012%
42	10.323	1370	1381	1388	rBV	428644	742236	80.03%	9.994%
43	10.390	1388	1392	1403	rVB	27795	47626	5.13%	0.641%
44	10.500	1406	1410	1416	rVB3	1977	3556	0.38%	0.048%
45	10.579	1416	1423	1430	rBV	74852	129742	13.99%	1.747%
46	10.628	1430	1431	1439	rVB3	2916	4062	0.44%	0.055%
47	10.701	1439	1443	1449	rBV	5338	8497	0.92%	0.114%
48	10.853	1464	1468	1470	rVB2	389	503	0.05%	0.007%
49	10.926	1473	1480	1494	rBV	119424	204266	22.02%	2.750%
50	11.073	1496	1504	1520	rBV	395703	642809	69.31%	8.655%
51	11.420	1554	1561	1565	rBV4	1346	2506	0.27%	0.034%
52	11.566	1579	1585	1588	rVV3	1269	2264	0.24%	0.030%
53	11.627	1588	1595	1604	rVV	371298	623146	67.19%	8.391%
54	11.707	1605	1608	1609	rVV2	1332	1496	0.16%	0.020%
55	11.731	1609	1612	1619	rVV3	1860	3256	0.35%	0.044%
56	11.841	1623	1630	1639	rVV4	10489	22853	2.46%	0.308%
57	11.944	1641	1647	1653	rVV2	7910	13847	1.49%	0.186%
58	12.018	1654	1659	1665	rVV3	1318	2760	0.30%	0.037%
59	12.097	1667	1672	1675	rVV4	389	890	0.10%	0.012%
60	12.170	1679	1684	1690	rVB3	1372	2368	0.26%	0.032%
61	12.237	1690	1695	1700	rBV2	1443	2176	0.23%	0.029%
62	12.335	1706	1711	1727	rVV	31904	52485	5.66%	0.707%
63	12.463	1728	1732	1737	rVV3	415	652	0.07%	0.009%
64	12.511	1737	1740	1744	rVB3	635	903	0.10%	0.012%
65	12.603	1747	1755	1761	rBV3	10527	19140	2.06%	0.258%
66	12.688	1763	1769	1777	rVV	148128	247662	26.70%	3.335%
67	12.755	1777	1780	1786	rVB5	669	1221	0.13%	0.016%
68	12.871	1794	1799	1800	rBV	357	479	0.05%	0.006%
69	12.932	1800	1809	1821	rVV4	4585	11748	1.27%	0.158%
70	13.048	1824	1828	1831	rVV3	2155	3145	0.34%	0.042%
71	13.097	1831	1836	1841	rVV2	7898	13479	1.45%	0.181%
72	13.152	1842	1845	1848	rVV4	2004	3305	0.36%	0.045%
73	13.206	1850	1854	1858	rVV5	1995	3822	0.41%	0.051%
74	13.280	1859	1866	1875	rVB6	3021	8947	0.96%	0.120%
75	13.395	1877	1885	1891	rBV2	21153	34649	3.74%	0.467%
76	13.475	1895	1898	1905	rVV6	1512	3680	0.40%	0.050%

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 Title : SFAM01.0

77	13.554	1905	1911	1922	rVV	297212	484766	52.27%	6.527%
78	13.645	1923	1926	1931	rVB3	4477	6728	0.73%	0.091%
79	13.712	1934	1937	1941	rVB3	406	375	0.04%	0.005%
80	13.767	1941	1946	1950	rBV5	926	1628	0.18%	0.022%
81	13.853	1950	1960	1972	rBV	279288	454715	49.03%	6.123%
82	14.017	1981	1987	1990	rBV5	2049	3294	0.36%	0.044%
83	14.066	1990	1995	2004	rVB	17847	29302	3.16%	0.395%
84	14.255	2024	2026	2030	rVB3	322	356	0.04%	0.005%
85	14.322	2030	2037	2046	rBV	17712	28148	3.03%	0.379%
86	14.420	2051	2053	2058	rVB3	293	359	0.04%	0.005%
87	14.517	2066	2069	2079	rVB3	1533	2883	0.31%	0.039%
88	14.615	2079	2085	2089	rBV3	1565	2668	0.29%	0.036%
89	14.718	2099	2102	2107	rVB2	1357	1929	0.21%	0.026%
90	14.877	2123	2128	2132	rBV4	837	1537	0.17%	0.021%
91	14.944	2135	2139	2145	rBV3	6328	10753	1.16%	0.145%
92	15.169	2171	2176	2185	rVB4	3424	6815	0.73%	0.092%
93	15.249	2185	2189	2194	rBV5	1014	1830	0.20%	0.025%
94	15.395	2204	2213	2220	rBV6	2050	4766	0.51%	0.064%
95	15.486	2223	2228	2236	rBV2	2689	5231	0.56%	0.070%
96	15.846	2282	2287	2297	rVB3	4335	7637	0.82%	0.103%
97	15.962	2299	2306	2309	rBV3	2394	4842	0.52%	0.065%
98	16.023	2314	2316	2319	rBV2	307	351	0.04%	0.005%
99	16.096	2324	2328	2332	rVB5	572	887	0.10%	0.012%
100	16.608	2409	2412	2415	rBV5	226	378	0.04%	0.005%

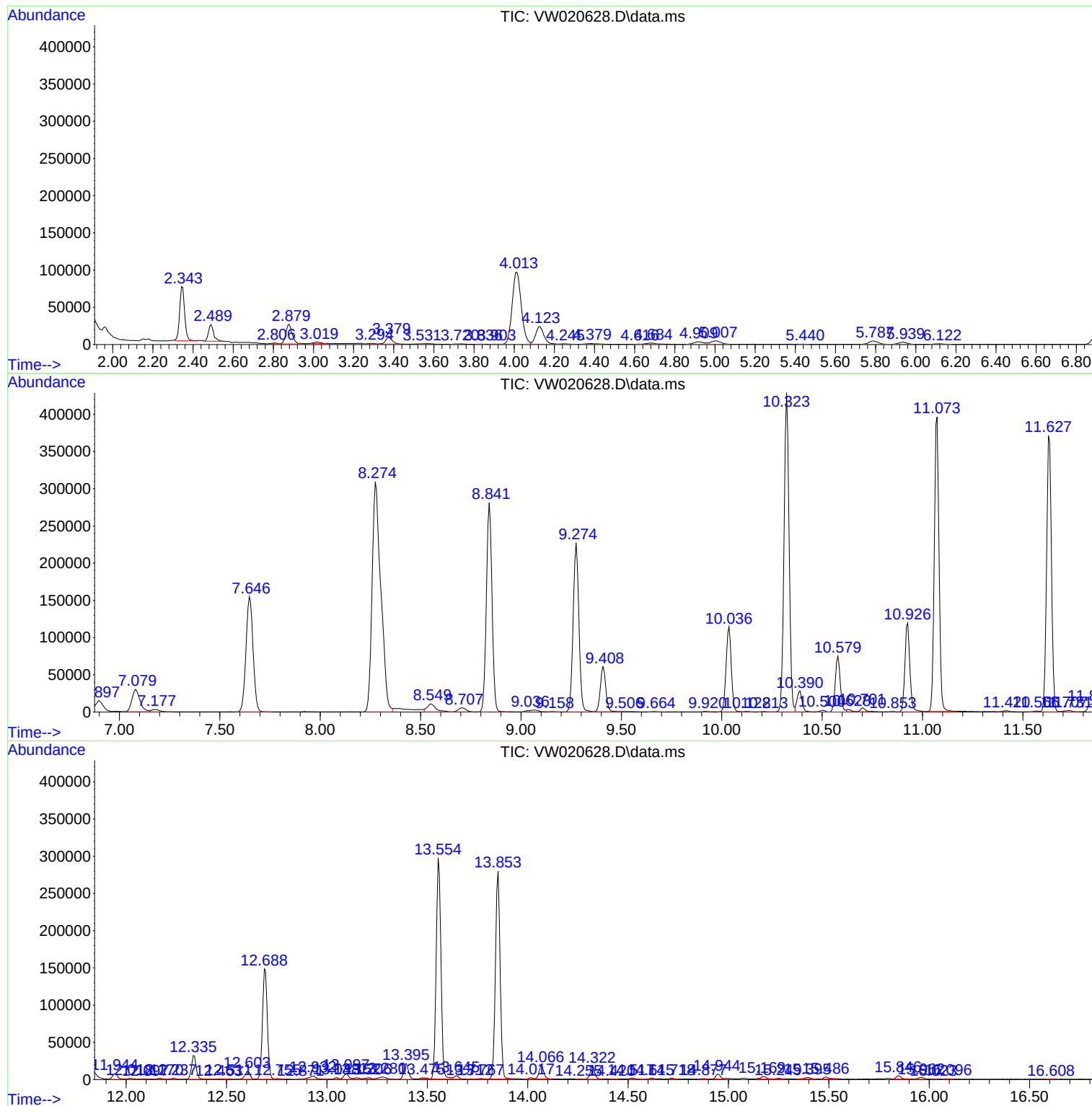
Sum of corrected areas: 7426788

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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



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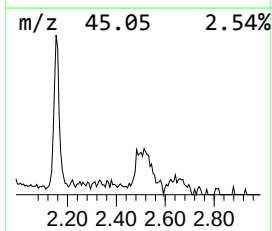
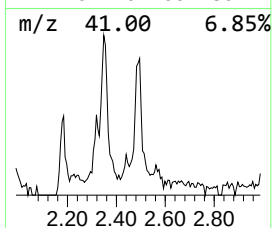
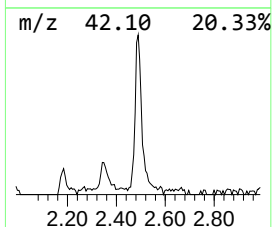
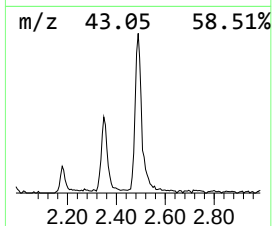
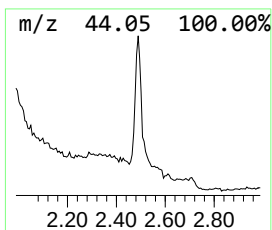
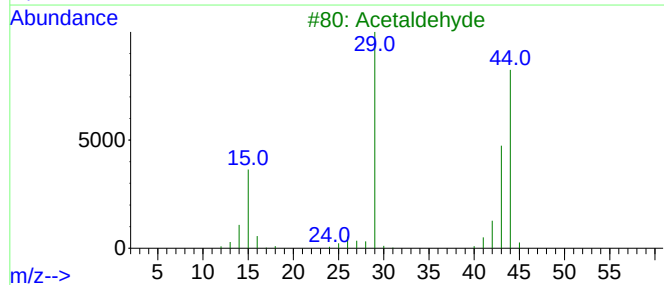
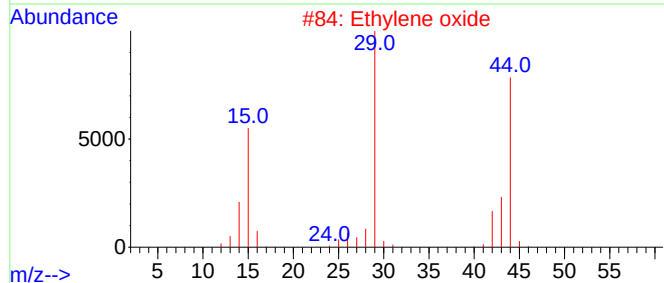
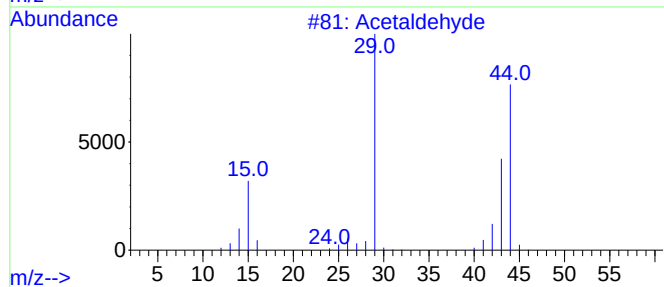
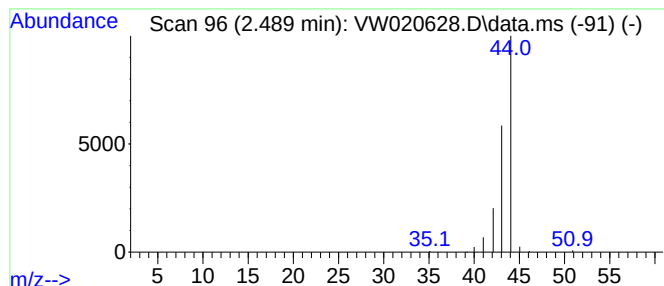
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.489	1.81 ug/L	39463	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetaldehyde	44	C2H4O	000075-07-0	9
2			Ethylene oxide	44	C2H4O	000075-21-8	9
3			Acetaldehyde	44	C2H4O	000075-07-0	9
4			Propane	44	C3H8	000074-98-6	5
5			Ethylene oxide	44	C2H4O	000075-21-8	5



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Instrument :
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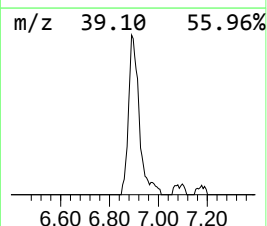
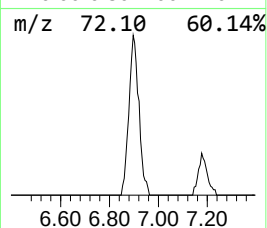
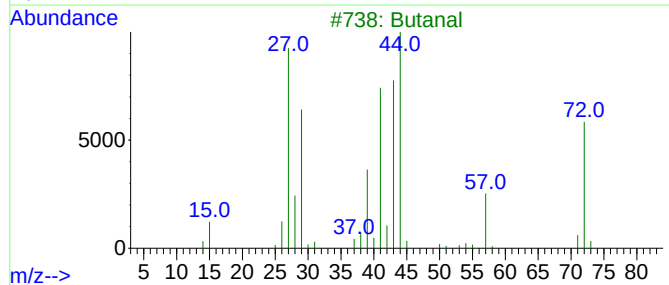
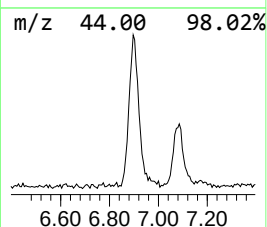
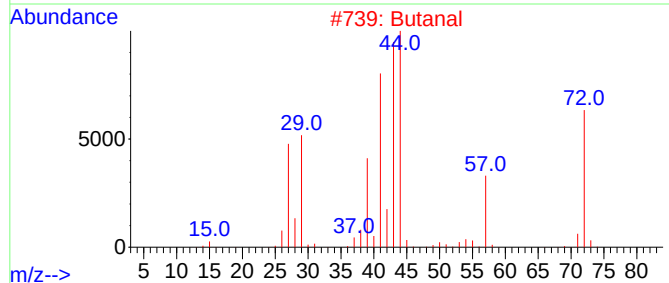
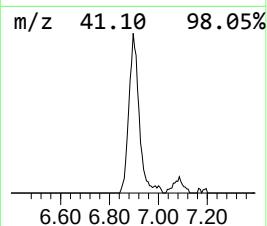
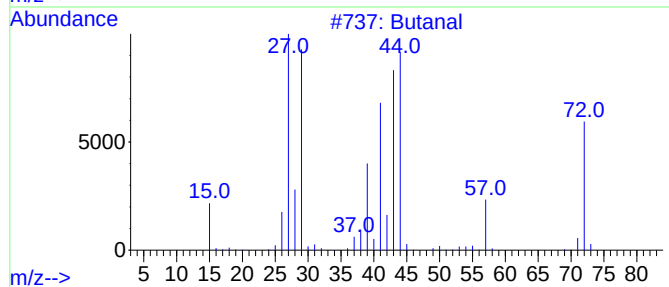
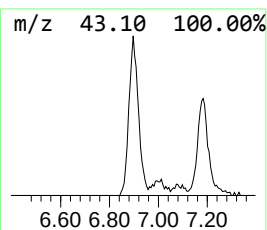
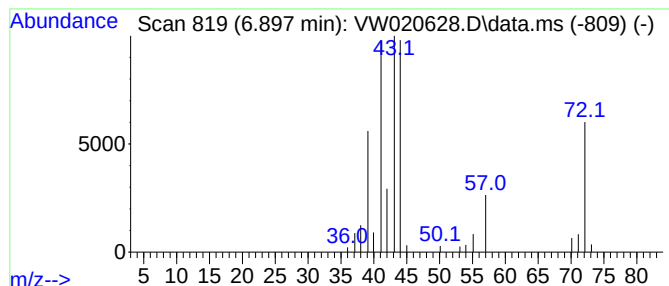
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 Butanal Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.897	1.95 ug/L	42421	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanal			72	C4H8O	000123-72-8	91
2	Butanal			72	C4H8O	000123-72-8	91
3	Butanal			72	C4H8O	000123-72-8	74
4	Butanal			72	C4H8O	000123-72-8	58
5	Butanal			72	C4H8O	000123-72-8	50



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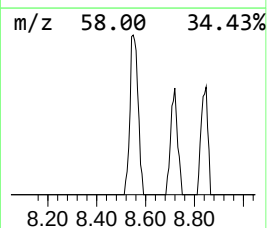
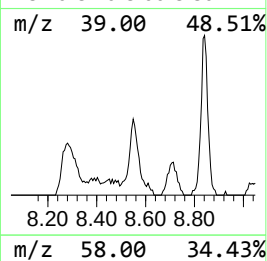
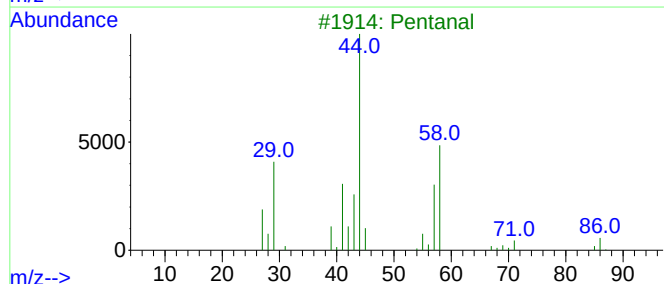
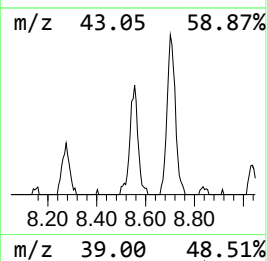
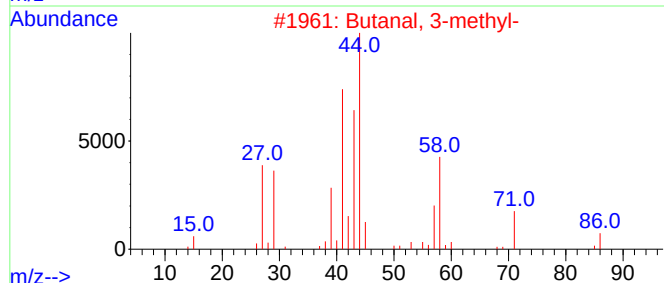
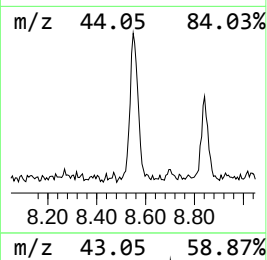
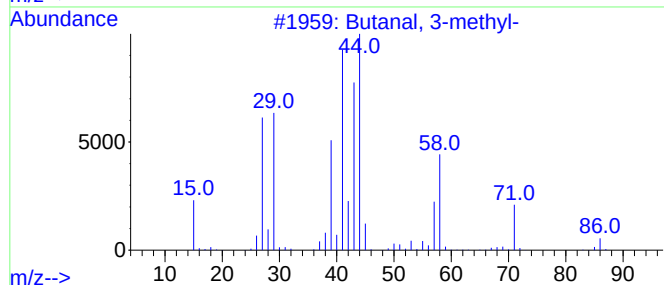
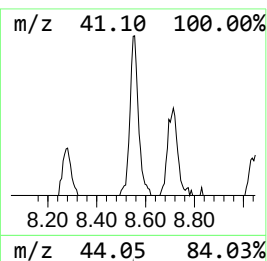
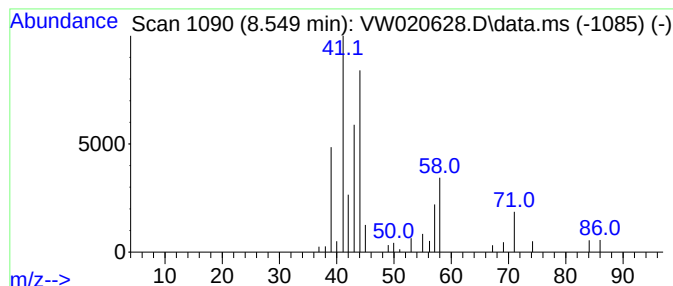
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Butanal, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.549	1.47 ug/L	31952	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanal, 3-methyl-	86	C5H10O	000590-86-3	58
2			Butanal, 3-methyl-	86	C5H10O	000590-86-3	52
3			Pentanal	86	C5H10O	000110-62-3	43
4			Butanal, 3-methyl-	86	C5H10O	000590-86-3	25
5			Butanal, 3-methyl-	86	C5H10O	000590-86-3	20



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Operator : SY/VA
Sample : M4263-11
Misc : 6.10g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GB7F6

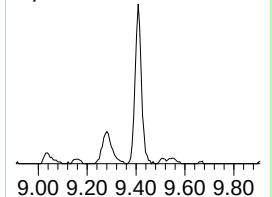
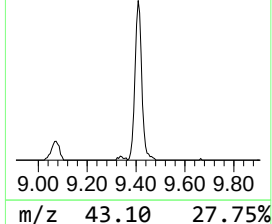
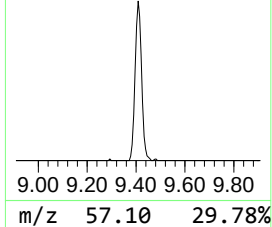
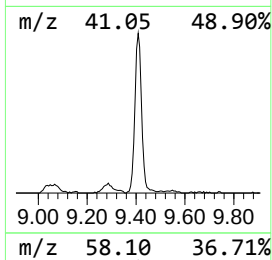
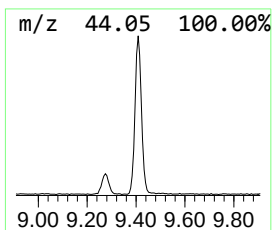
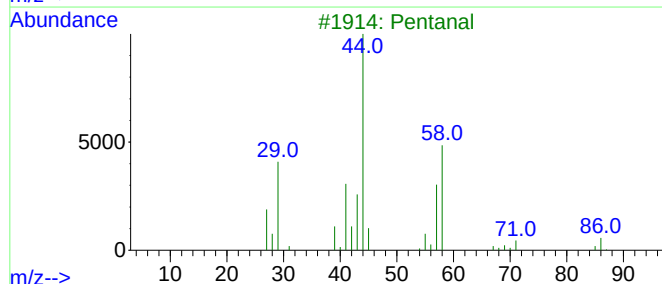
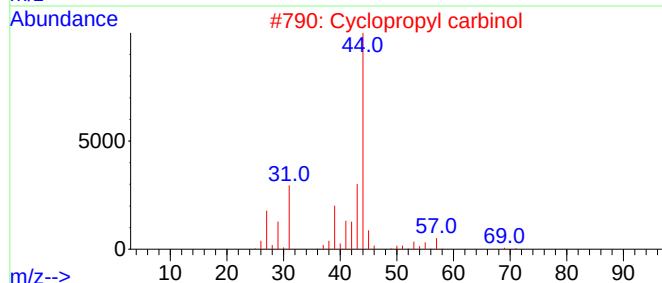
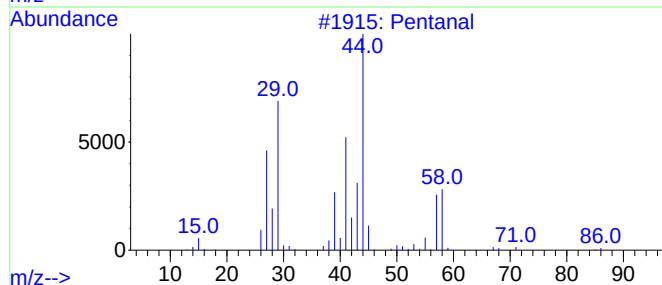
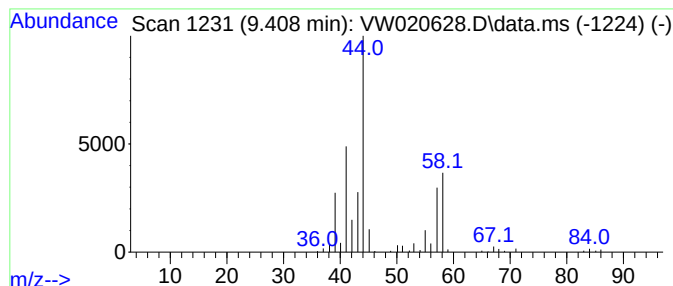
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Quant Title : SFAM01.0

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

Peak Number 4 Pentanal Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.408	5.19 ug/L	113106	1,4-Difluorobenzene	8.841

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanal	86	C5H10O	000110-62-3	86
2			Cyclopropyl carbinol	72	C4H8O	002516-33-8	59
3			Pentanal	86	C5H10O	000110-62-3	56
4			Pentanal	86	C5H10O	000110-62-3	49
5			Butanal, 3-methyl-	86	C5H10O	000590-86-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101921\
Data File : VW020628.D
Acq On : 20 Oct 2021 00:25
Operator : SY/VA
Sample : M4263-11
Misc : 6.10g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GB7F6

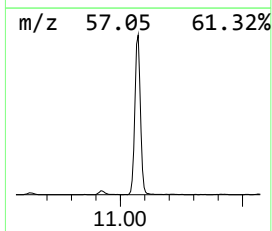
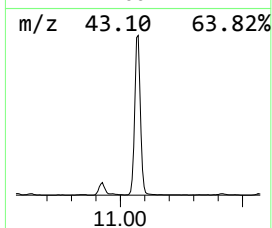
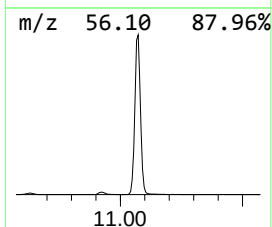
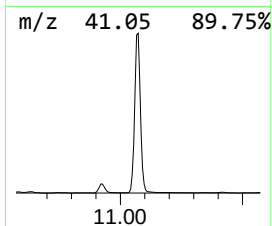
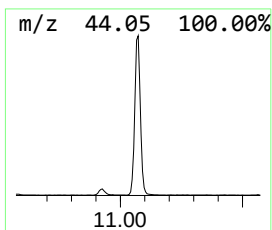
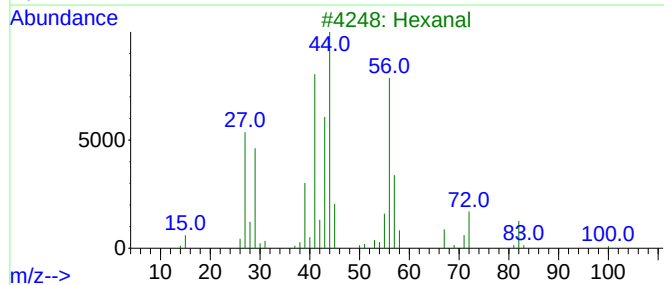
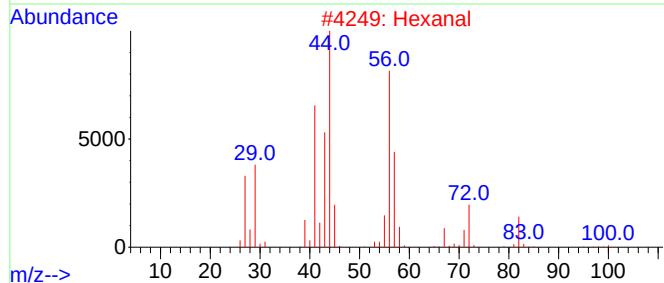
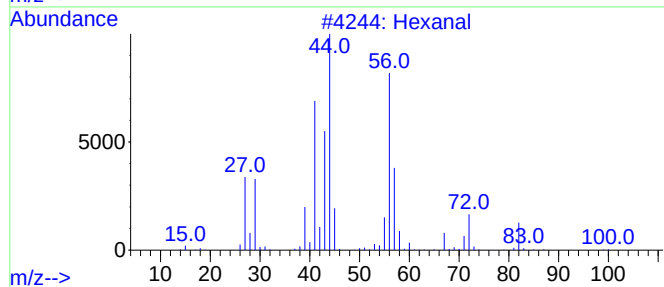
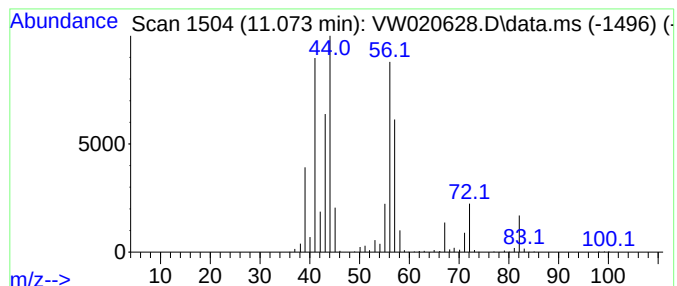
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Quant Title : SFAM01.0

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

Peak Number 6 Hexanal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.073	25.79 ug/L	642809	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal			100	C6H12O	000066-25-1	87
2	Hexanal			100	C6H12O	000066-25-1	80
3	Hexanal			100	C6H12O	000066-25-1	80
4	Hexanal			100	C6H12O	000066-25-1	72
5	Hexanal			100	C6H12O	000066-25-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101921\
Data File : VW020628.D
Acq On : 20 Oct 2021 00:25
Operator : SY/VA
Sample : M4263-11
Misc : 6.10g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GB7F6

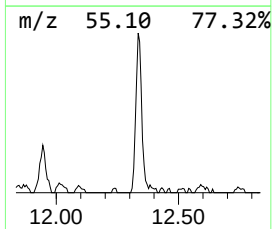
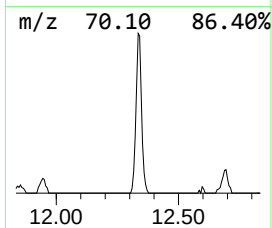
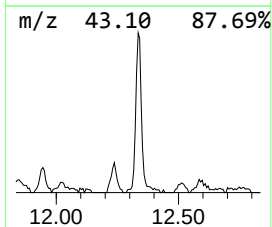
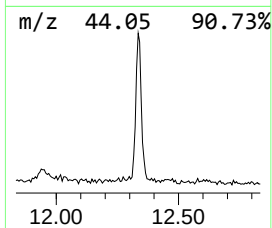
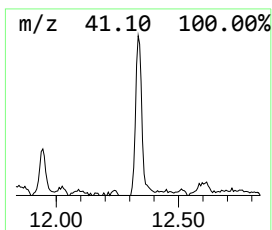
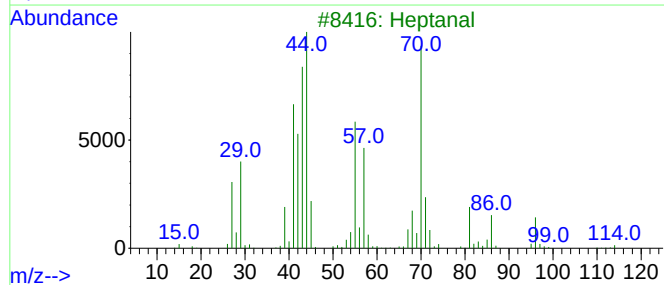
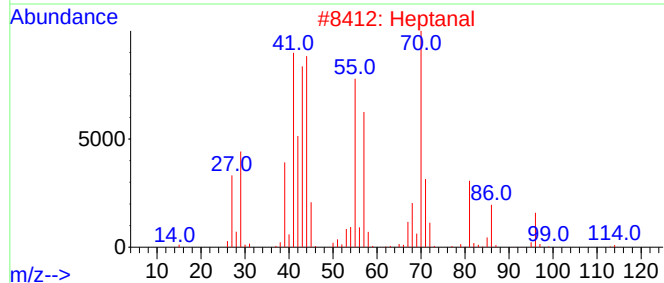
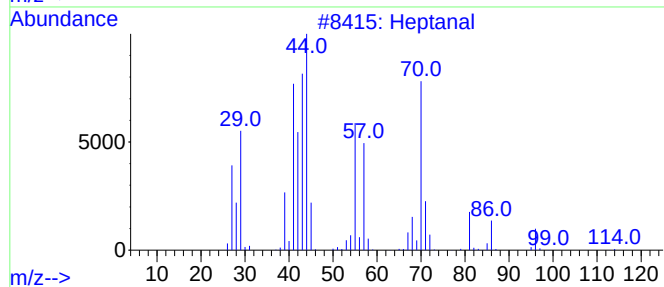
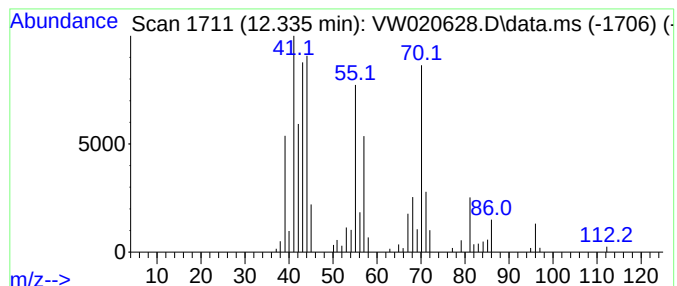
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Quant Title : SFAM01.0

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

Peak Number 7 Heptanal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.335	2.11 ug/L	52485	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanal			114	C7H14O	000111-71-7	90
2	Heptanal			114	C7H14O	000111-71-7	90
3	Heptanal			114	C7H14O	000111-71-7	90
4	Heptanal			114	C7H14O	000111-71-7	78
5	Heptanal			114	C7H14O	000111-71-7	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101921\
Data File : VW020628.D
Acq On : 20 Oct 2021 00:25
Operator : SY/VA
Sample : M4263-11
Misc : 6.10g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GB7F6

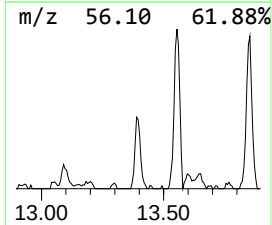
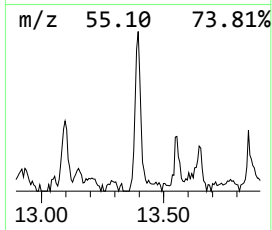
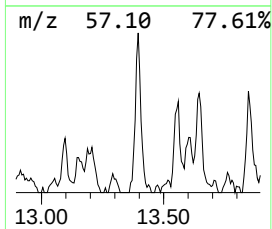
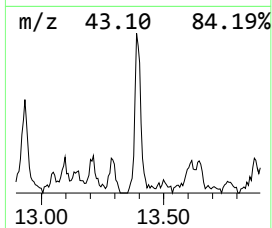
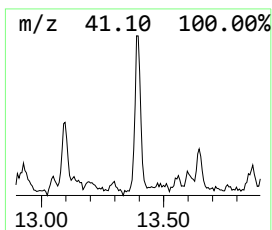
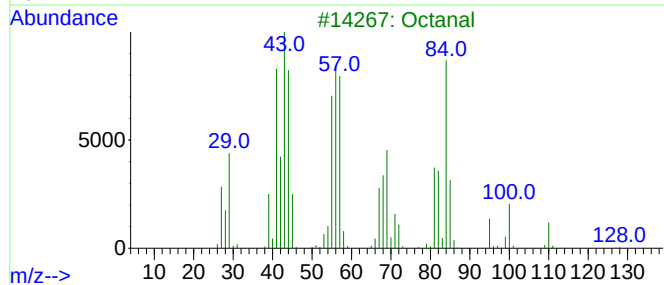
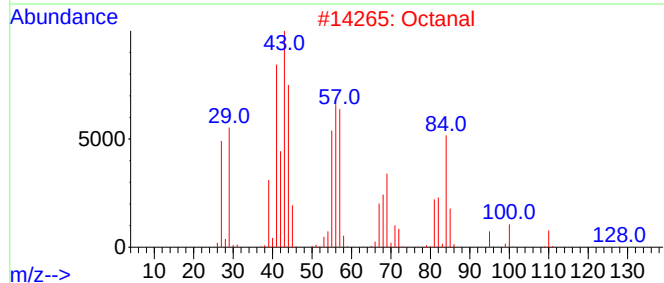
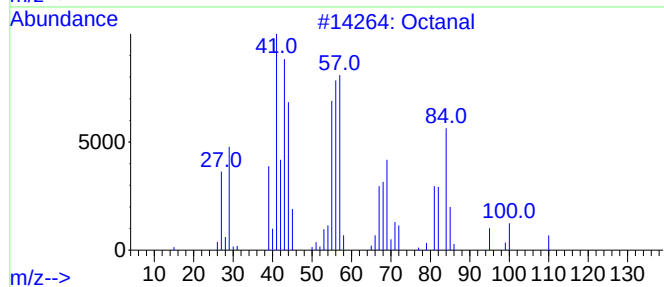
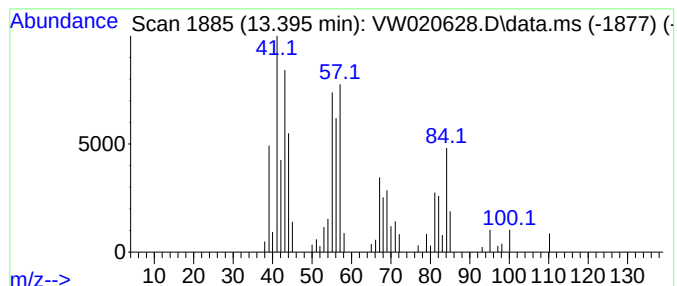
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Quant Title : SFAM01.0

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

Peak Number 8 Octanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.395	1.79 ug/L	34649	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanal			128	C8H16O	000124-13-0	91
2	Octanal			128	C8H16O	000124-13-0	86
3	Octanal			128	C8H16O	000124-13-0	80
4	Octanal			128	C8H16O	000124-13-0	53
5	Pentane, 3-methylene-			84	C6H12	000760-21-4	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101921\
Data File : VW020628.D
Acq On : 20 Oct 2021 00:25
Operator : SY/VA
Sample : M4263-11
Misc : 6.10g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GB7F6

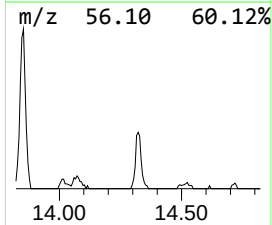
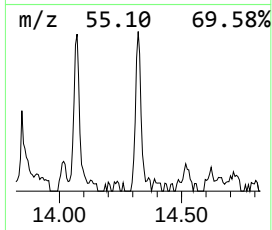
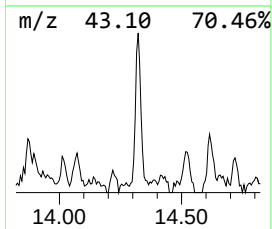
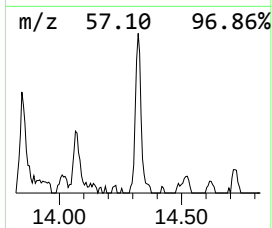
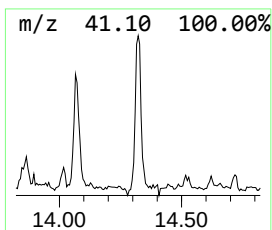
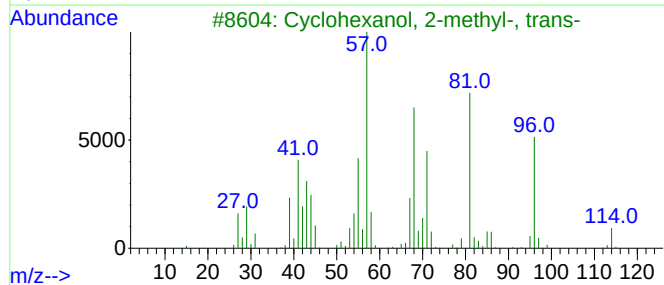
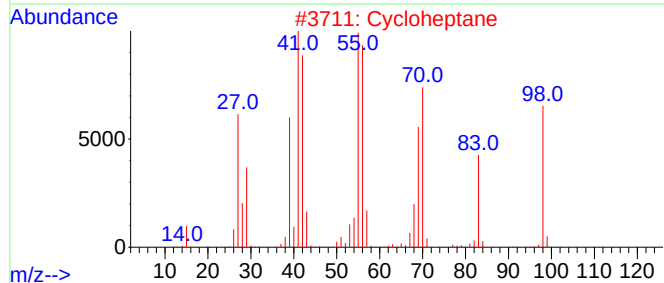
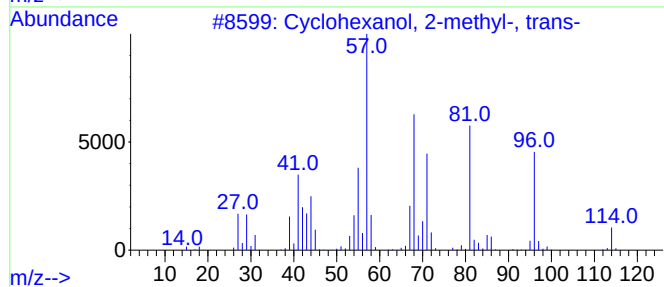
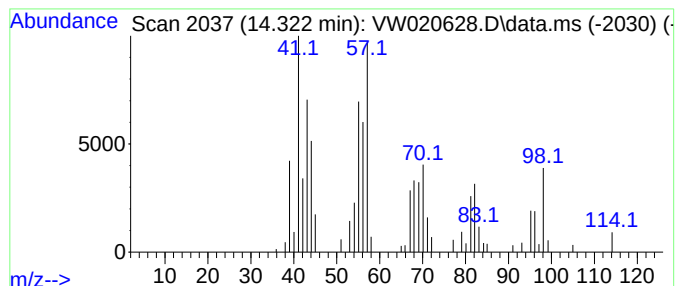
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Quant Title : SFAM01.0

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

Peak Number 10 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.322	1.45 ug/L	28148	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexanol, 2-methyl-, trans-	114	C7H14O	007443-52-9	38
2		Cycloheptane	98	C7H14	000291-64-5	38
3		Cyclohexanol, 2-methyl-, trans-	114	C7H14O	007443-52-9	38
4		2-Nonen-1-ol, (E)-	142	C9H18O	031502-14-4	35
5		5-Methyl-2-hexene,c&t	98	C7H14	003404-62-4	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101921\
Data File : VW020628.D
Acq On : 20 Oct 2021 00:25
Operator : SY/VA
Sample : M4263-11
Misc : 6.10g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
GB7F6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Butanal	6.897	2.0	ug/L	42421	1	8.841	545230	25.0
Butanal, 3-methyl-	8.549	1.5	ug/L	31952	1	8.841	545230	25.0
Pentanal	9.408	5.2	ug/L	113106	1	8.841	545230	25.0
Hexanal	11.073	25.8	ug/L	642809	2	11.634	623146	25.0
Heptanal	12.335	2.1	ug/L	52485	2	11.634	623146	25.0
Octanal	13.395	1.8	ug/L	34649	3	13.560	484766	25.0
unknown-02	14.322	1.5	ug/L	28148	3	13.560	484766	25.0