

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
 Data File : VX018750.D
 Acq On : 02 Oct 2020 16:16
 Operator : JC/SP
 Sample : L4171-13DL2 50X
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3W0DL2

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.386	42	49	55	rBV	89665	97357	2.08%	0.197%
2	1.697	94	100	106	rBV	78976	110971	2.37%	0.224%
3	2.343	199	206	217	rVB	186571	285058	6.08%	0.575%
4	2.825	276	285	287	rBV2	115285	226885	4.84%	0.458%
5	2.867	287	292	305	rVB	761672	1569376	33.49%	3.168%
6	3.148	329	338	354	rBV	1694987	3829949	81.72%	7.731%
7	3.501	385	396	411	rBV	2132671	4686681	100.00%	9.461%
8	4.111	486	496	508	rVB3	46578	137454	2.93%	0.277%
9	4.276	511	523	531	rBV	31816	104790	2.24%	0.212%
10	4.373	531	539	551	rVB	101836	296958	6.34%	0.599%
11	4.562	557	570	584	rBV	60306	265228	5.66%	0.535%
12	5.141	654	665	681	rBV2	109341	328808	7.02%	0.664%
13	6.044	799	813	828	rBV2	238869	724904	15.47%	1.463%
14	6.830	933	942	952	rBV	179599	436103	9.31%	0.880%
15	7.367	1022	1030	1039	rBV	145075	317658	6.78%	0.641%
16	8.372	1189	1195	1202	rBV	100838	168864	3.60%	0.341%
17	8.696	1237	1248	1259	rBV	400967	632929	13.50%	1.278%
18	8.994	1289	1297	1306	rVB2	72671	117017	2.50%	0.236%
19	9.427	1362	1368	1378	rBV	464426	683057	14.57%	1.379%
20	10.098	1472	1478	1488	rBV	382938	531410	11.34%	1.073%
21	11.000	1621	1626	1633	rBV	55901	75047	1.60%	0.151%
22	11.232	1659	1664	1672	rVB	159960	213051	4.55%	0.430%
23	11.341	1676	1682	1689	rBV	988910	1251393	26.70%	2.526%
24	11.421	1689	1695	1702	rVV	1427387	2739796	58.46%	5.531%
25	11.494	1702	1707	1714	rVB	1317613	1682363	35.90%	3.396%
26	11.652	1727	1733	1741	rBV	280389	342685	7.31%	0.692%
27	11.793	1751	1756	1765	rVB	2052752	2501376	53.37%	5.049%
28	11.902	1765	1774	1776	rBV	121984	157646	3.36%	0.318%
29	11.927	1776	1778	1783	rVB	158417	185471	3.96%	0.374%
30	12.006	1786	1791	1795	rVV	337812	424472	9.06%	0.857%
31	12.061	1795	1800	1805	rVB2	489498	703844	15.02%	1.421%
32	12.128	1805	1811	1816	rVB	161224	199354	4.25%	0.402%
33	12.213	1820	1825	1831	rVB	33073	48504	1.03%	0.098%
34	12.286	1831	1837	1838	rBV	670147	727662	15.53%	1.469%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
 Data File : VX018750.D
 Acq On : 02 Oct 2020 16:16
 Operator : JC/SP
 Sample : L4171-13DL2 50X
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3W0DL2

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.1
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
 Title : TRACE VOA SOM01.0

35	12.305	1838	1840	1844	rVV	1004420	1194354	25.48%	2.411%
36	12.353	1844	1848	1858	rVB2	2269123	3309354	70.61%	6.680%
37	12.445	1858	1863	1868	rBV2	124975	171526	3.66%	0.346%
38	12.500	1868	1872	1878	rVB	503121	610350	13.02%	1.232%
39	12.573	1878	1884	1886	rBV	1367417	1928742	41.15%	3.893%
40	12.591	1886	1887	1892	rVB	1258062	1196199	25.52%	2.415%
41	12.652	1892	1897	1901	rBV	2560487	2921586	62.34%	5.898%
42	12.683	1901	1902	1907	rVB	50751	47197	1.01%	0.095%
43	12.750	1907	1913	1921	rBV3	224479	406767	8.68%	0.821%
44	12.835	1922	1927	1931	rBV3	45693	57482	1.23%	0.116%
45	12.890	1931	1936	1940	rVB	629166	768565	16.40%	1.551%
46	12.963	1940	1948	1951	rBV	1636900	2021334	43.13%	4.080%
47	13.000	1951	1954	1960	rVB	2555337	3023810	64.52%	6.104%
48	13.061	1960	1964	1966	rBV	90786	107056	2.28%	0.216%
49	13.109	1970	1972	1975	rVB	63313	53164	1.13%	0.107%
50	13.207	1982	1988	1992	rBV	488603	604576	12.90%	1.220%
51	13.250	1992	1995	1998	rVV	44800	49105	1.05%	0.099%
52	13.292	1998	2002	2004	rVV2	125678	181541	3.87%	0.366%
53	13.329	2004	2008	2017	rVV2	1254849	1814544	38.72%	3.663%
54	13.408	2017	2021	2027	rVV	153599	174212	3.72%	0.352%
55	13.542	2038	2043	2047	rBV2	38266	62281	1.33%	0.126%
56	13.628	2051	2057	2064	rVV2	814340	1246169	26.59%	2.516%
57	13.719	2067	2072	2076	rVB	129645	168860	3.60%	0.341%
58	13.969	2107	2113	2115	rBV2	59538	87211	1.86%	0.176%
59	13.999	2115	2118	2124	rVB3	153049	210510	4.49%	0.425%
60	14.115	2133	2137	2143	rBV	79387	98188	2.10%	0.198%
61	14.256	2156	2160	2166	rBV	49445	76843	1.64%	0.155%
62	14.359	2172	2177	2182	rBV	47428	64274	1.37%	0.130%
63	14.676	2222	2229	2234	rBV	51835	78567	1.68%	0.159%

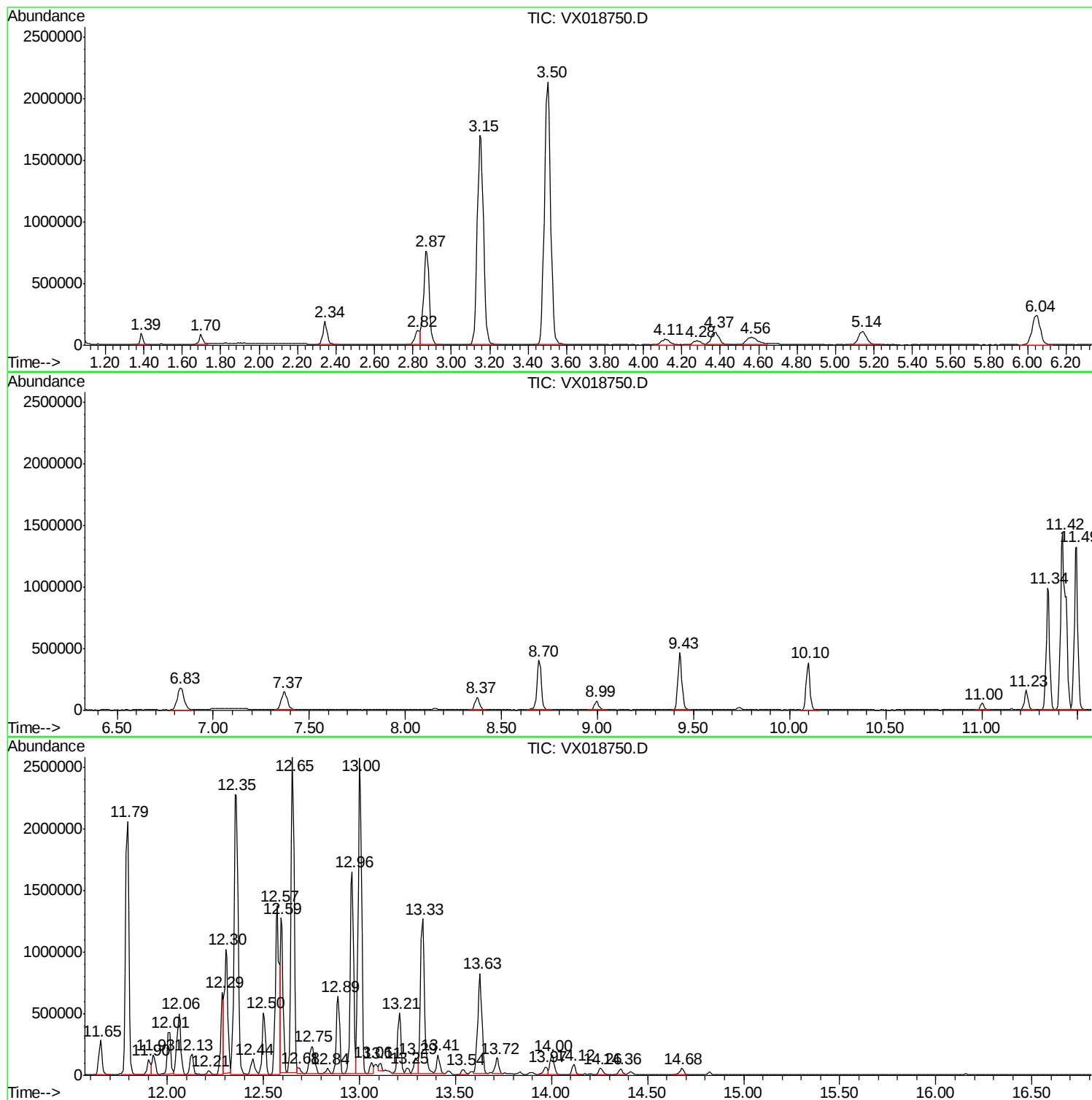
Sum of corrected areas: 49538488

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

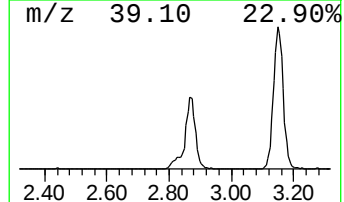
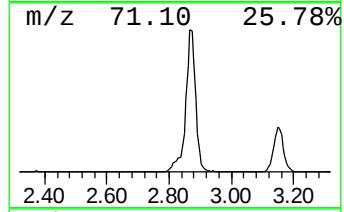
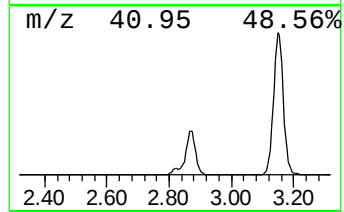
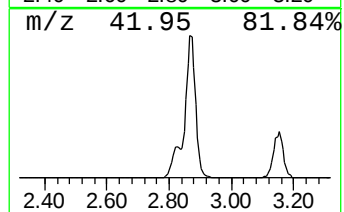
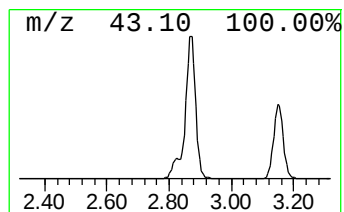
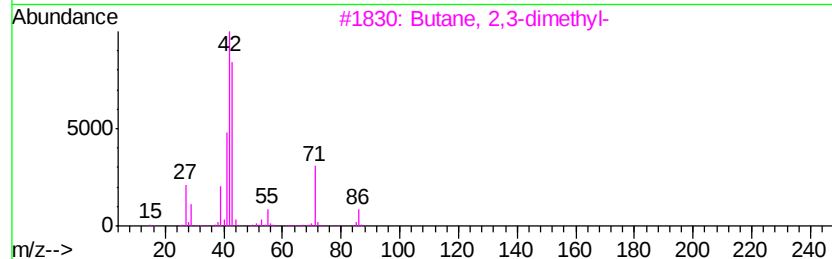
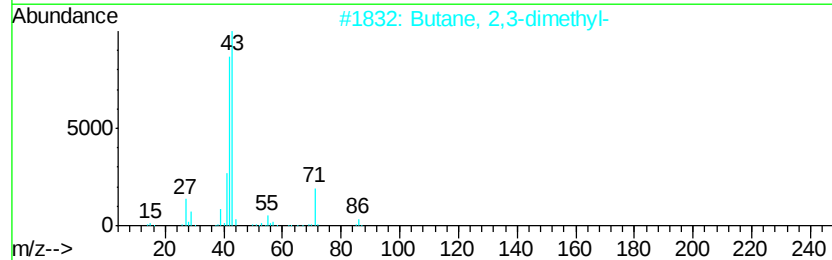
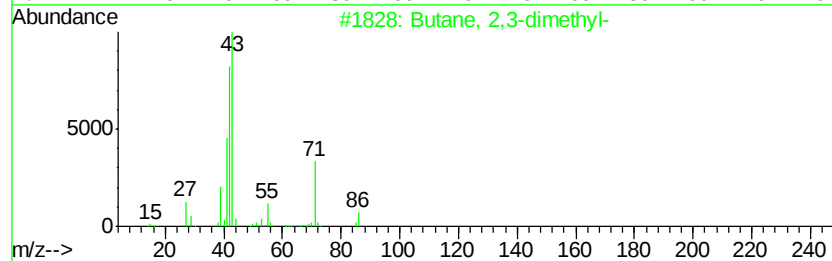
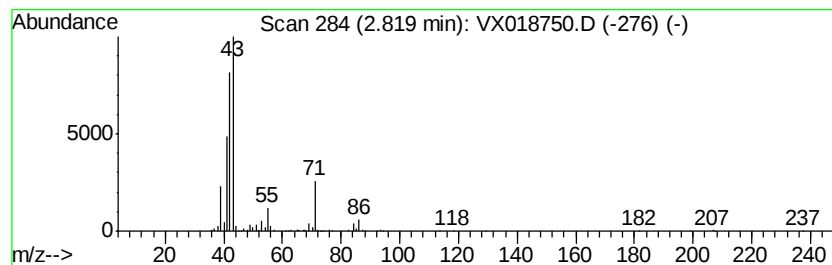
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.82	2.60 ug/L	226885	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	83
2		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
3		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
4		Methylamine, N-(1-propylbutylide...	127	C8H17N	010599-78-7	33
5		Pentane, 2-methyl-	86	C6H14	000107-83-5	28



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

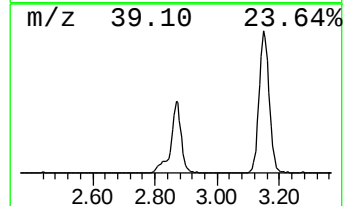
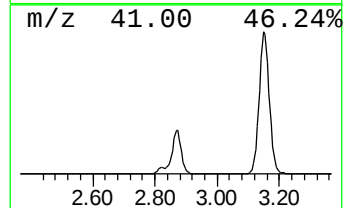
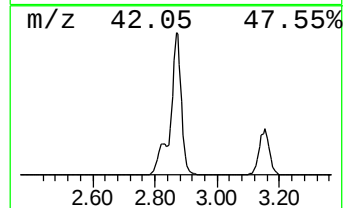
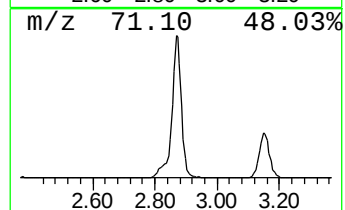
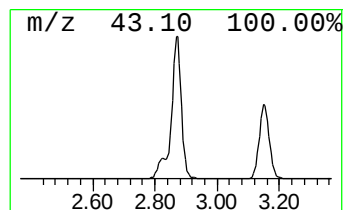
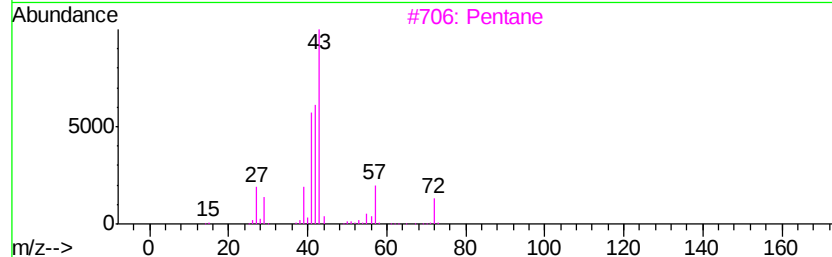
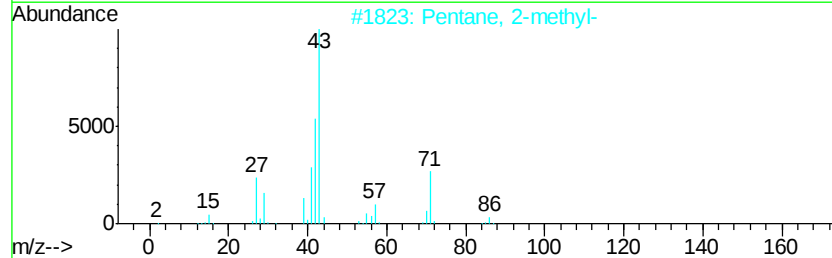
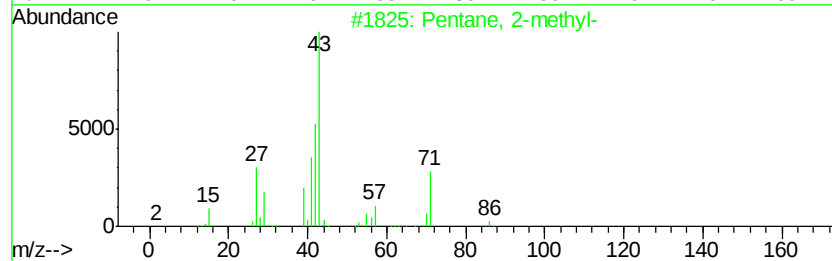
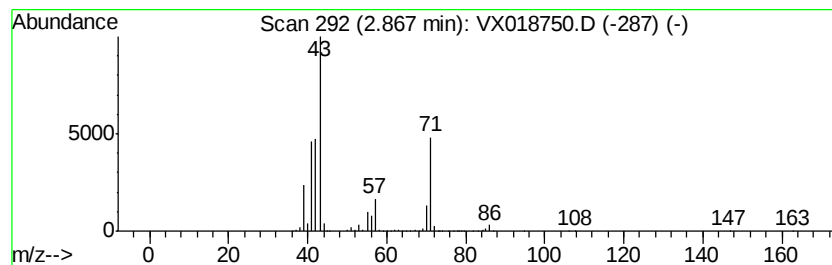
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	17.99 ug/L	1569380	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	83
2		Pentane, 2-methyl-	86	C6H14	000107-83-5	83
3		Pentane	72	C5H12	000109-66-0	47
4		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
5		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

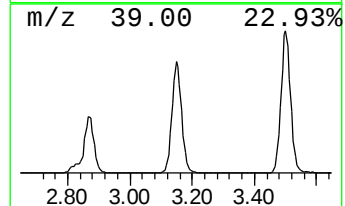
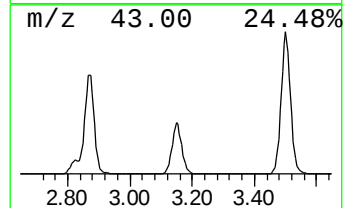
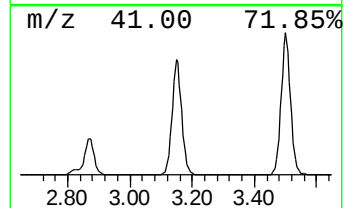
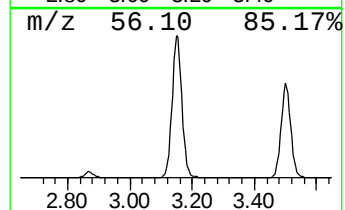
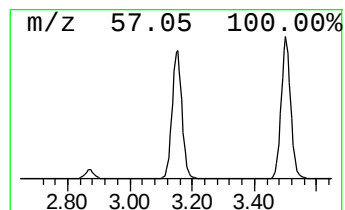
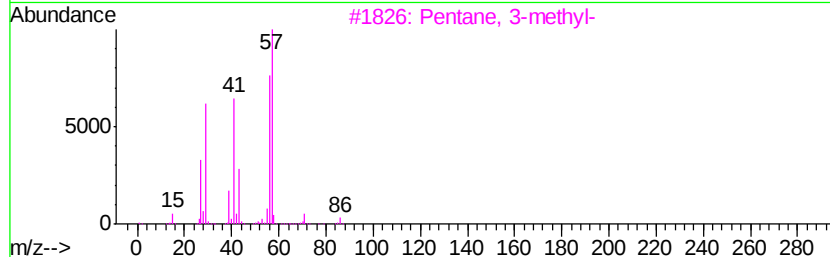
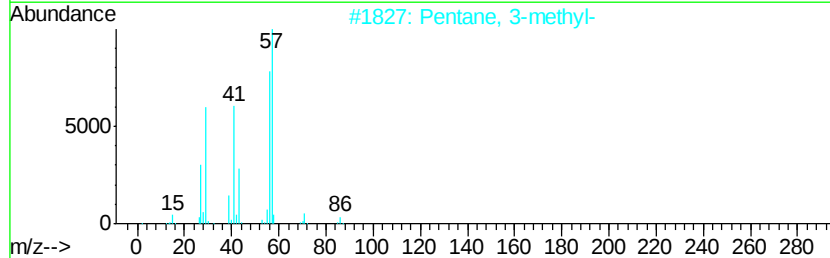
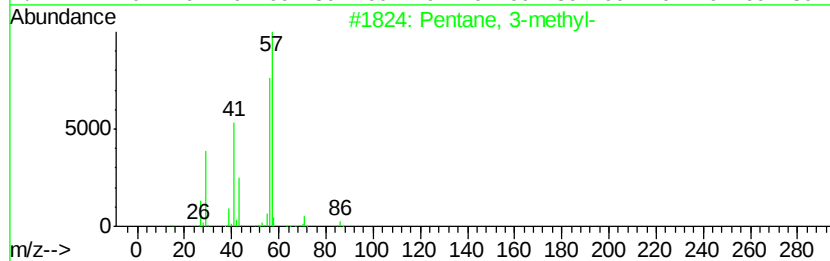
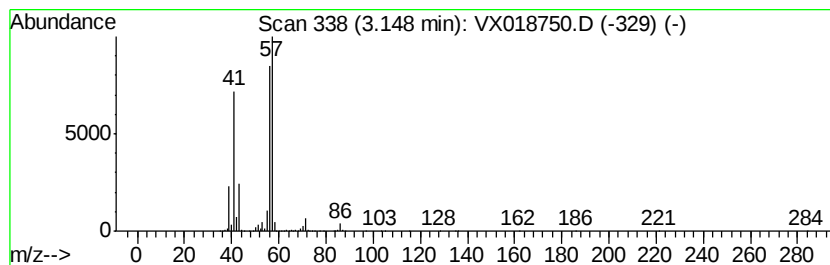
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	43.91 ug/L	3829950	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	90
3		Pentane, 3-methyl-	86	C6H14	000096-14-0	90
4		3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
5		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

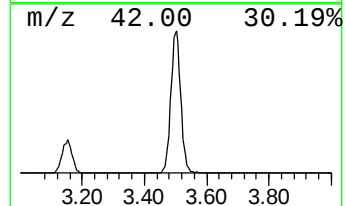
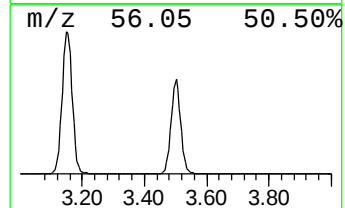
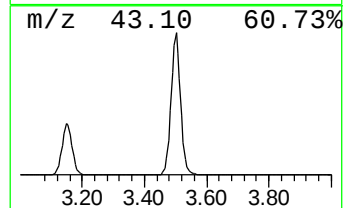
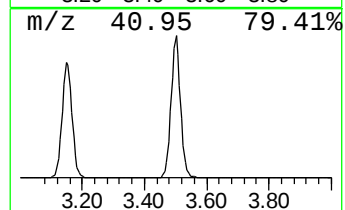
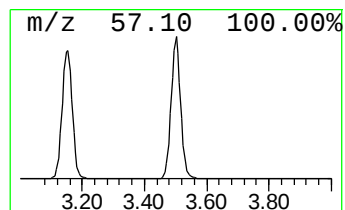
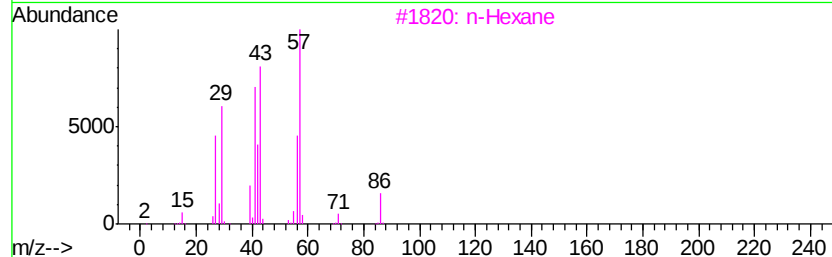
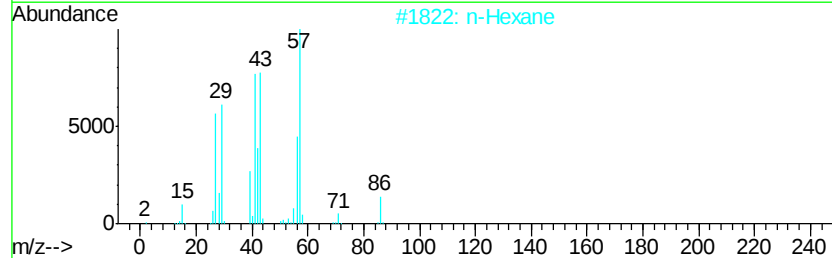
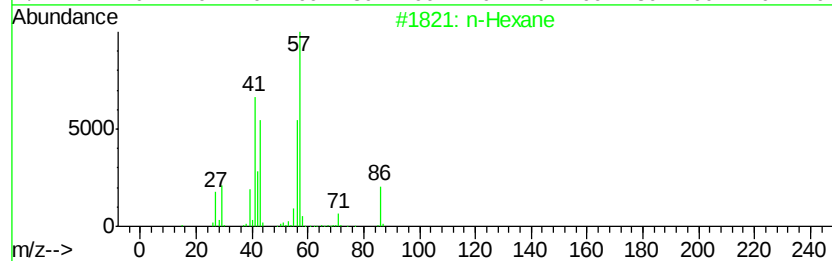
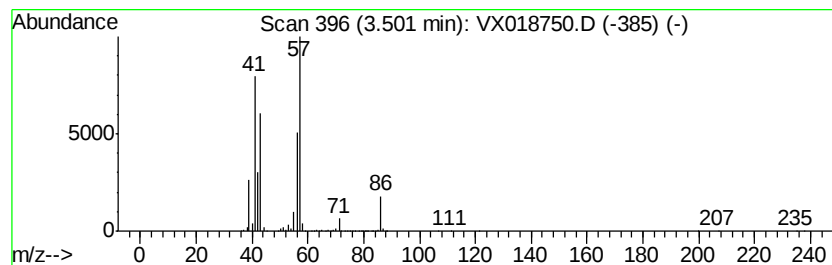
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	53.73 ug/L	4686680	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	94
2		n-Hexane	86	C6H14	000110-54-3	72
3		n-Hexane	86	C6H14	000110-54-3	59
4		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	50
5		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	38



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

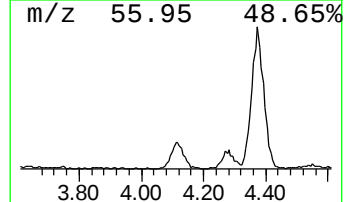
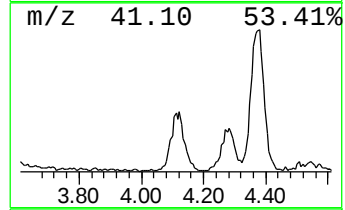
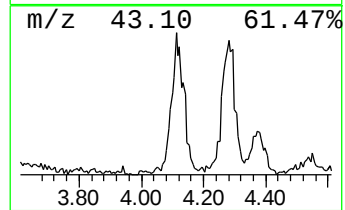
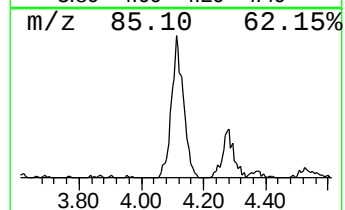
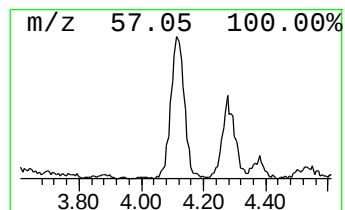
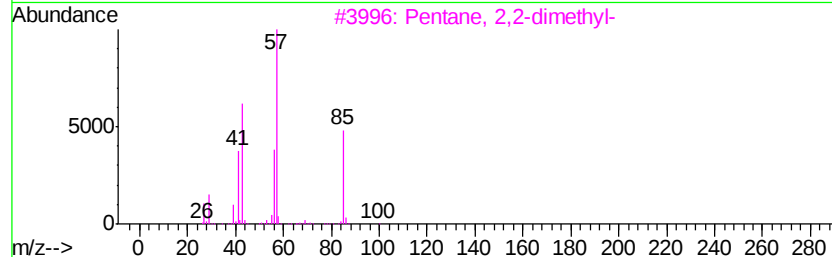
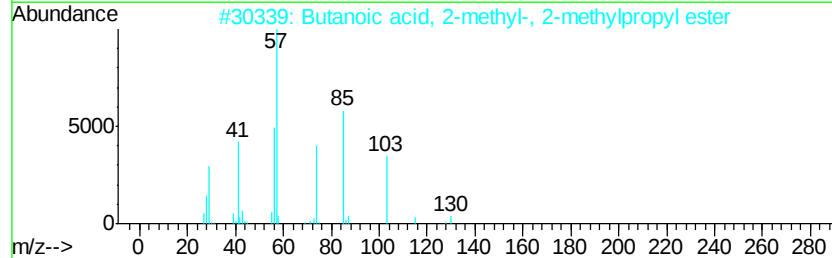
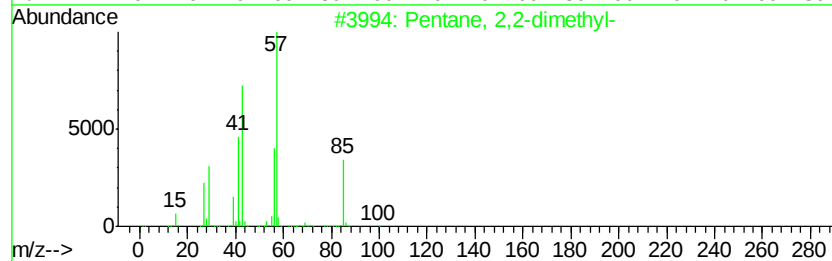
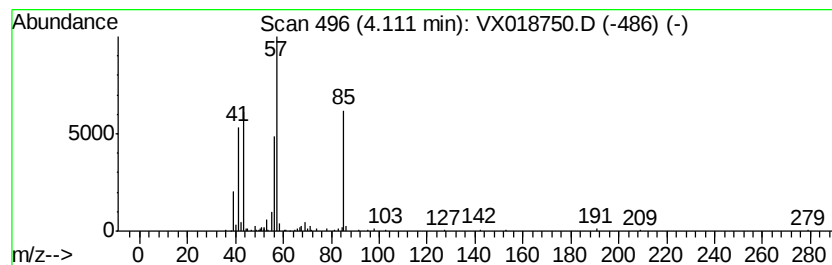
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.11	1.58 ug/L	137454	1,4-Difluorobenzene	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	72
2	Butanoic acid, 2-methyl-, 2-meth...	158	C9H18O2	002445-67-2	72
3	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64
4	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	53
5	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

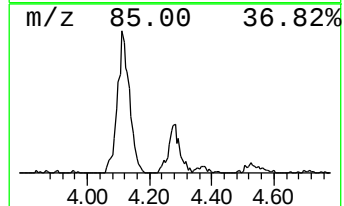
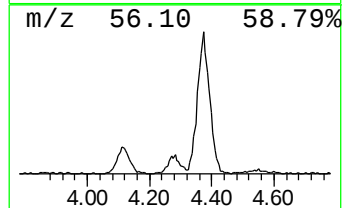
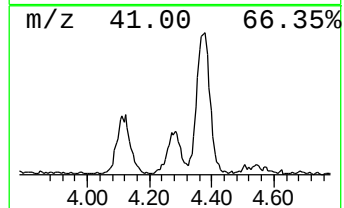
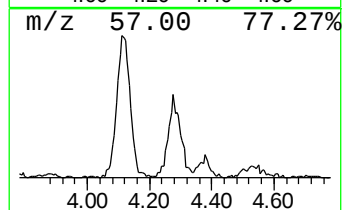
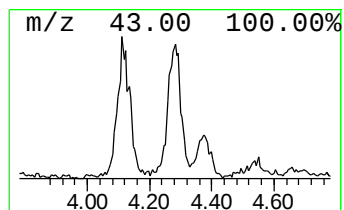
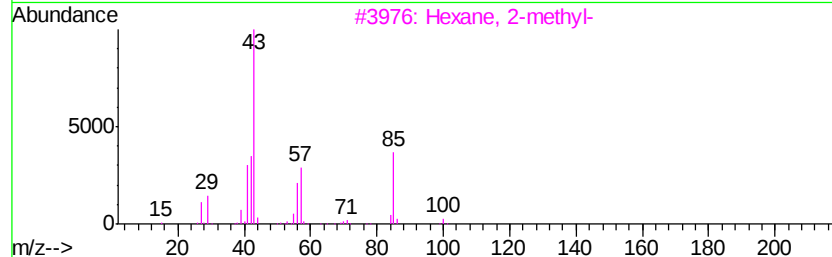
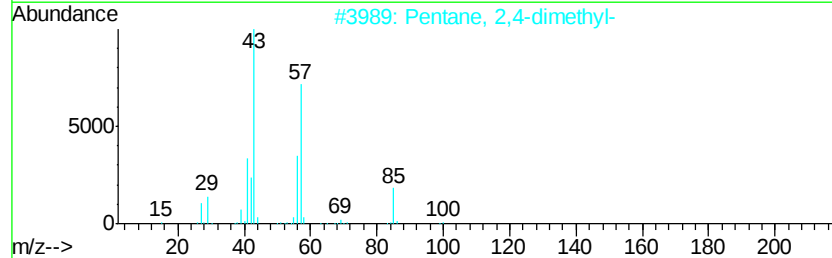
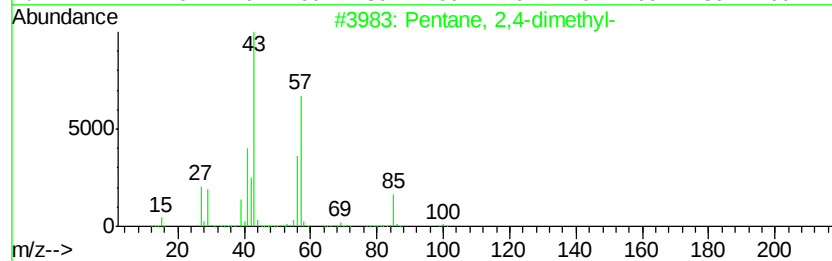
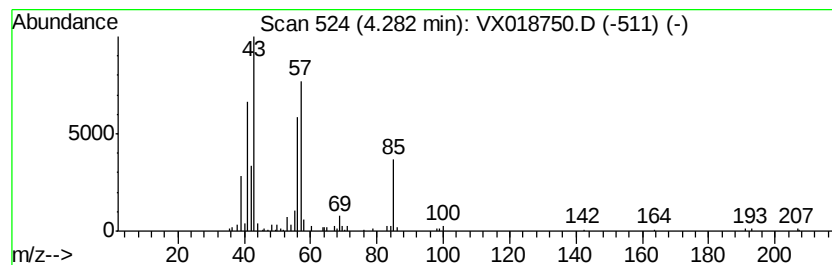
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	1.20 ug/L	104790	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	78
2		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	78
3		Hexane, 2-methyl-	100	C7H16	000591-76-4	64
4		Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	59
5		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

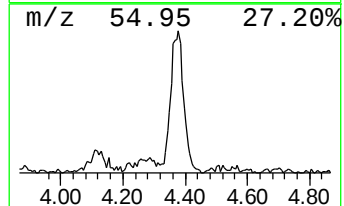
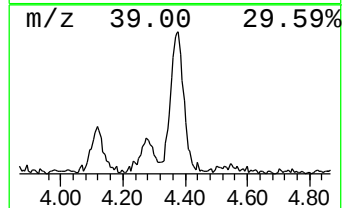
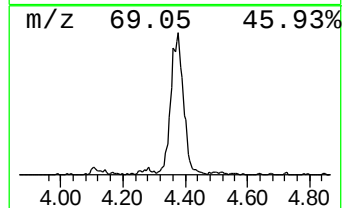
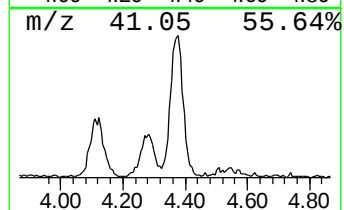
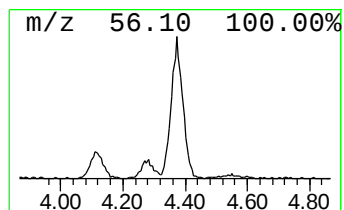
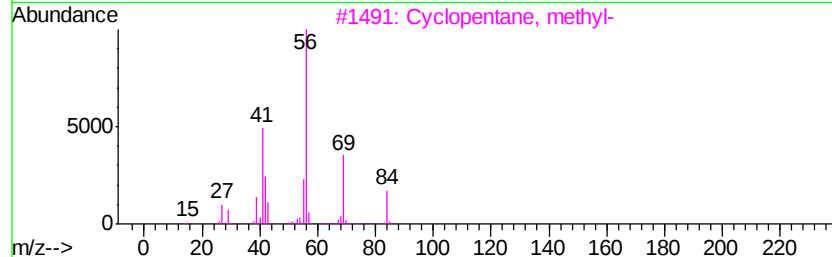
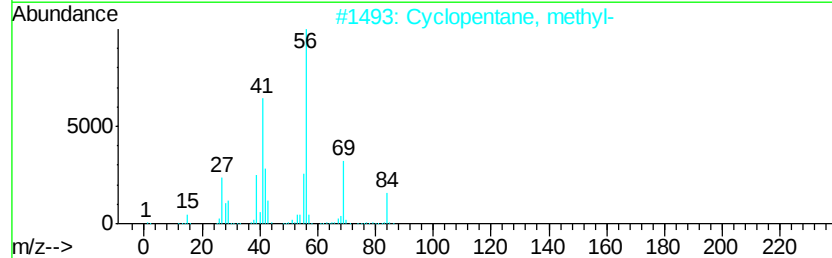
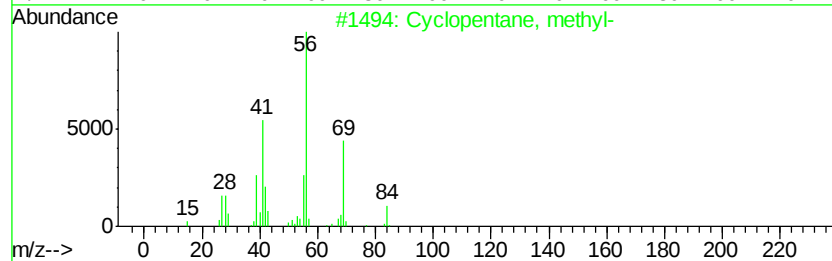
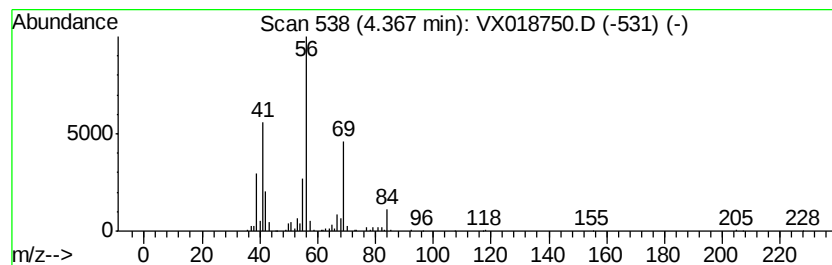
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Cyclic4.37 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	3.40 ug/L	296958	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	80
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	72
4		Pentane, 2-cyclopropyl-	112	C8H16	005458-16-2	47
5		2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

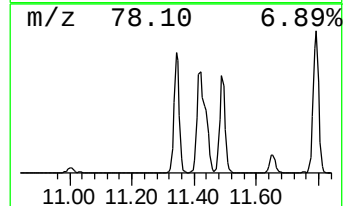
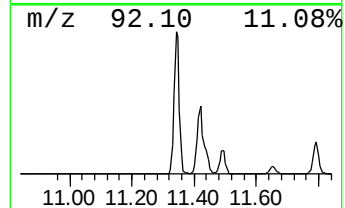
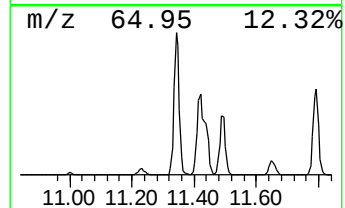
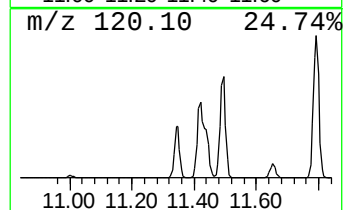
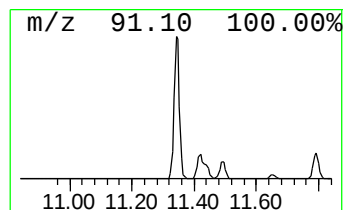
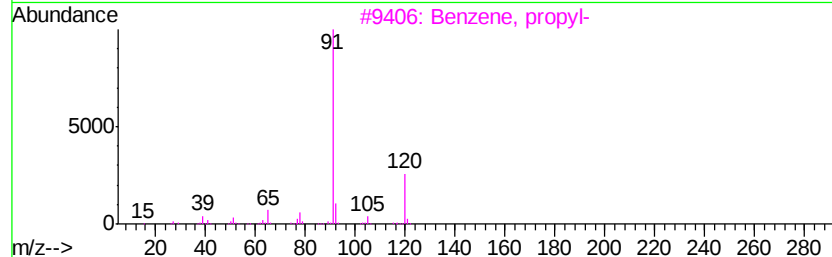
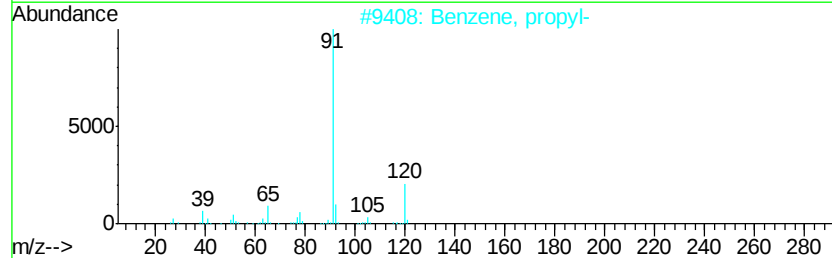
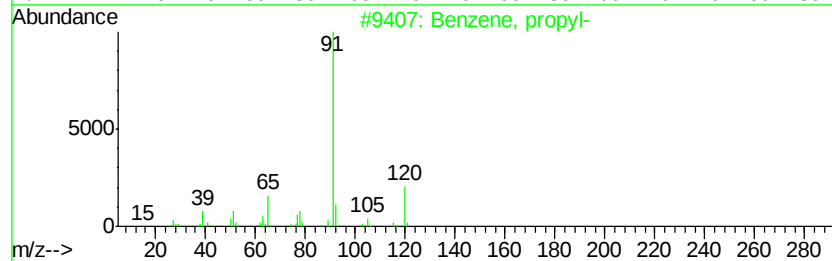
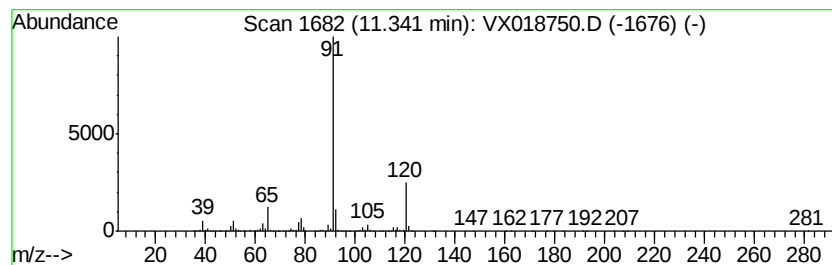
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.34	8.89 ug/L	1251390	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	90
2		Benzene, propyl-	120	C9H12	000103-65-1	90
3		Benzene, propyl-	120	C9H12	000103-65-1	83
4		Propanenitrile, 3-[(phenylmethyl)...	160	C10H12N2	000706-03-6	72
5		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

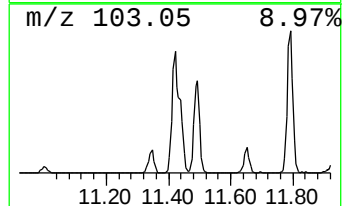
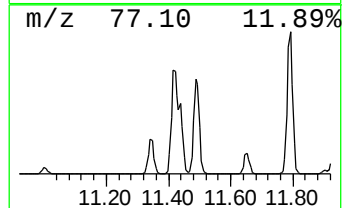
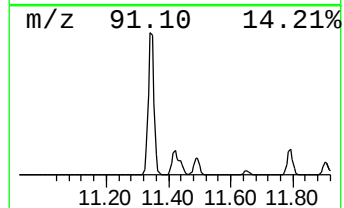
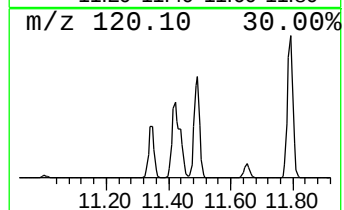
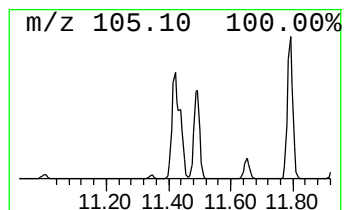
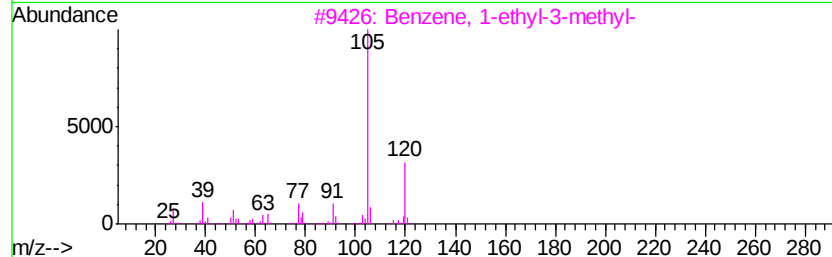
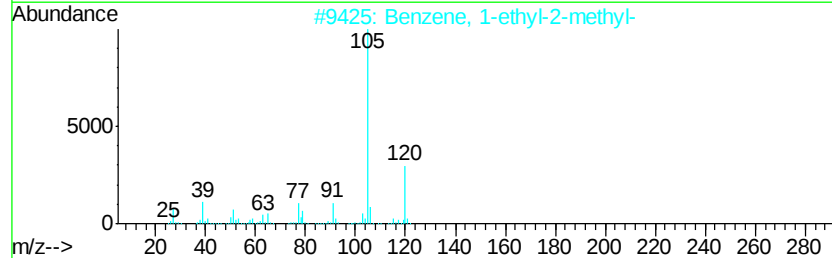
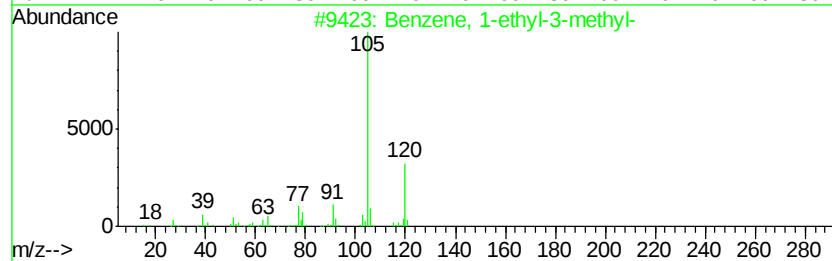
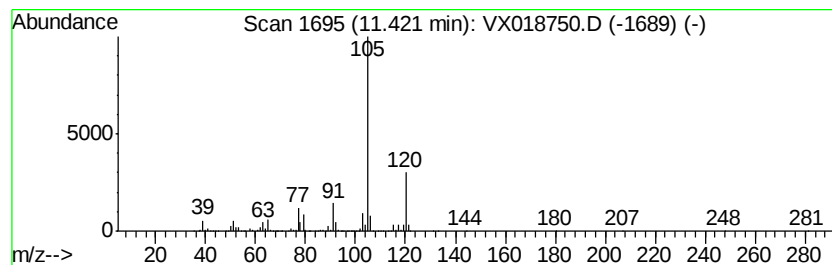
Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.42	19.46 ug/L	2739800	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

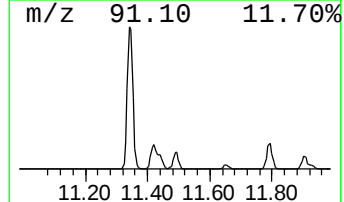
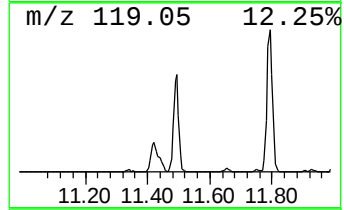
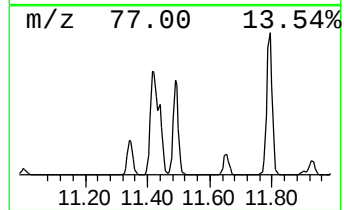
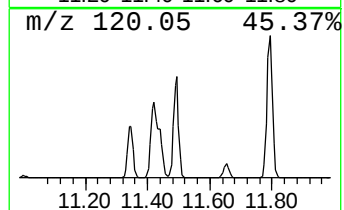
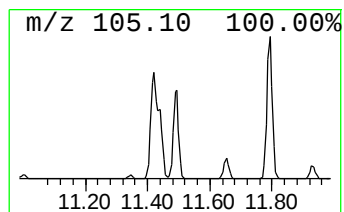
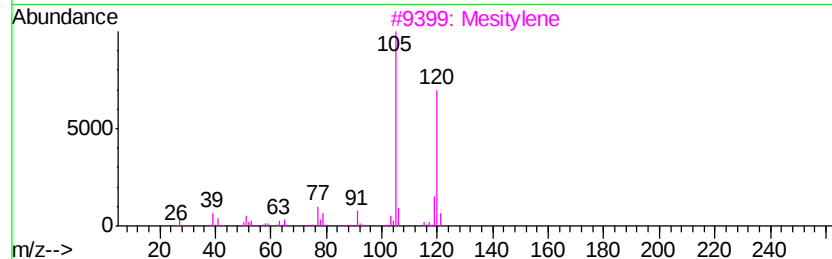
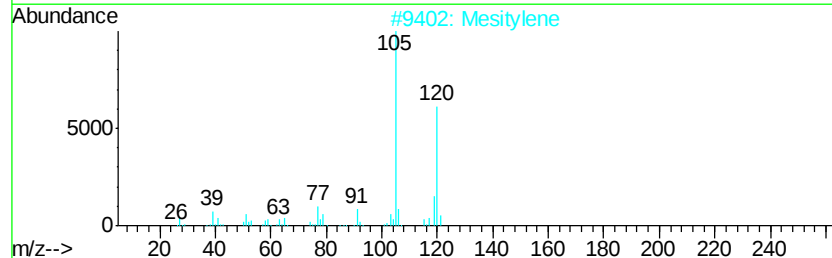
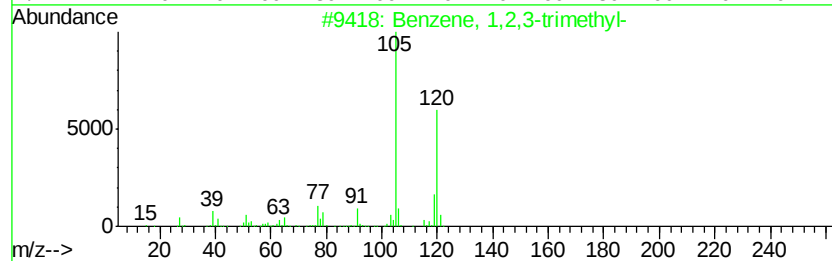
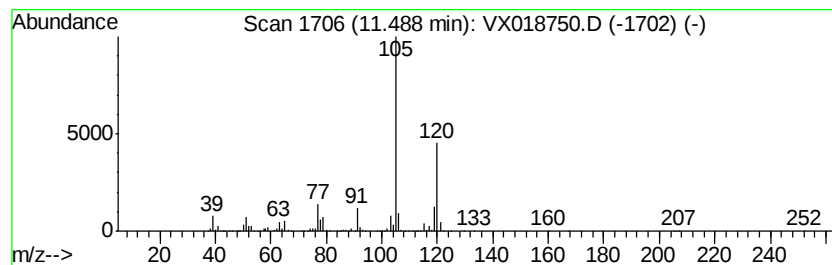
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.49	11.95 ug/L	1682360	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Mesitylene	120	C9H12	000108-67-8	97
3		Mesitylene	120	C9H12	000108-67-8	94
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5		Mesitylene	120	C9H12	000108-67-8	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

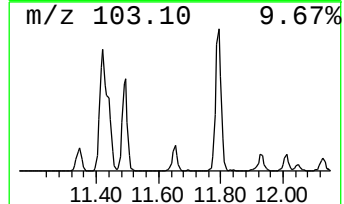
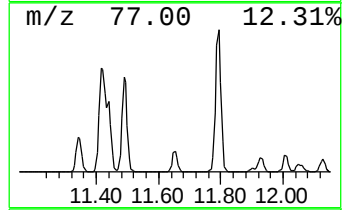
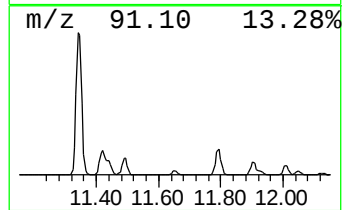
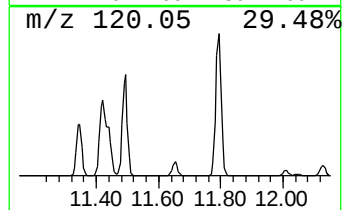
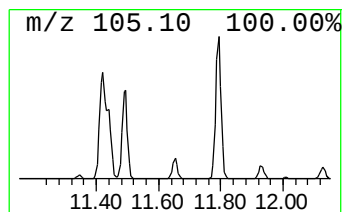
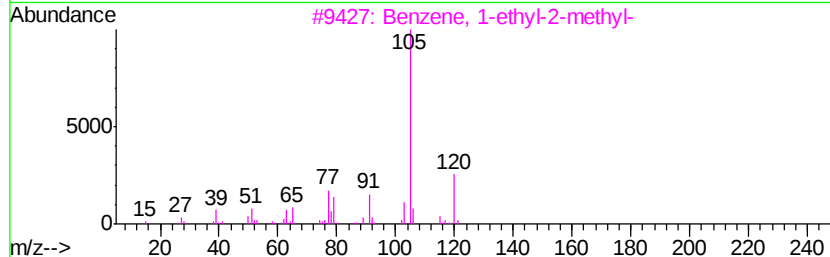
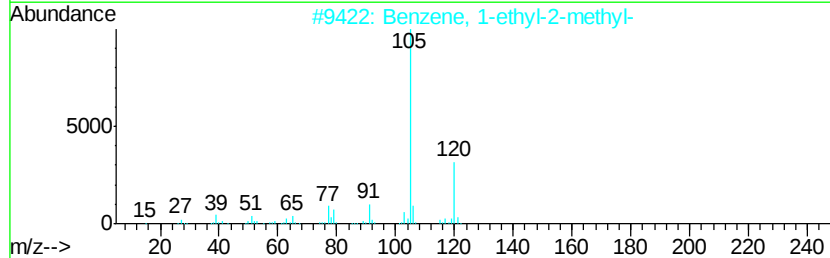
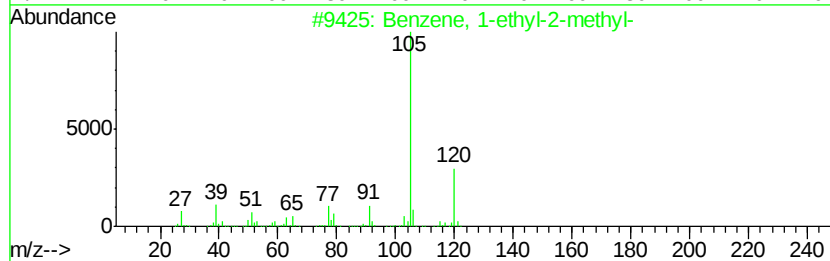
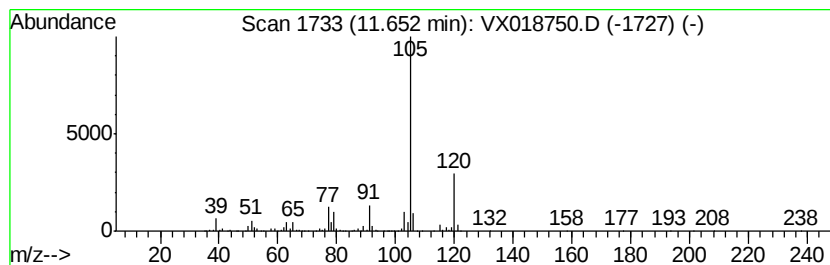
Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Benzene, 1-ethyl-2-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	2.43 ug/L	342685	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3	94
2	Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3	94
3	Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3	93
4	Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

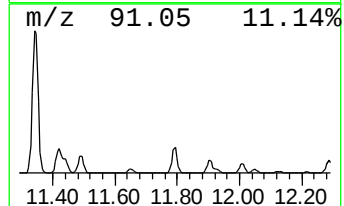
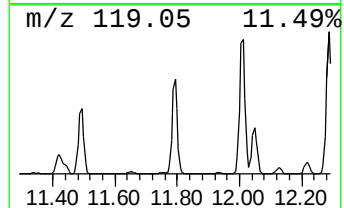
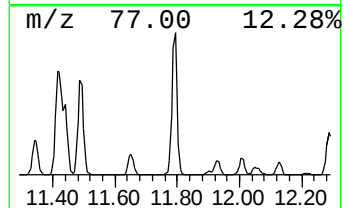
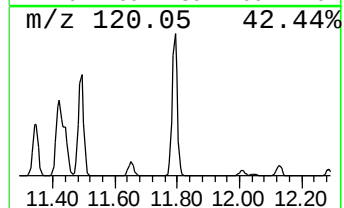
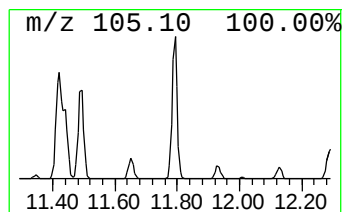
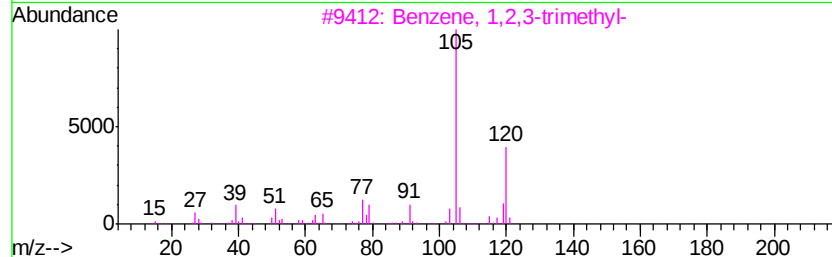
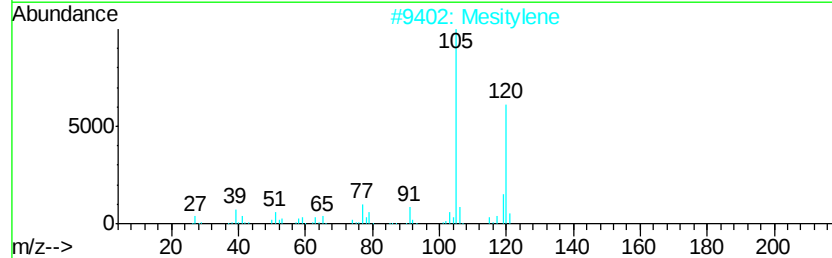
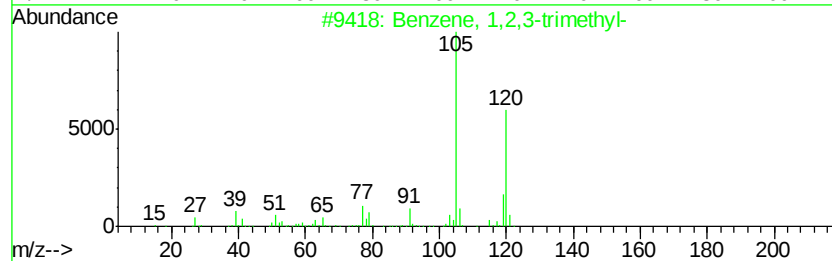
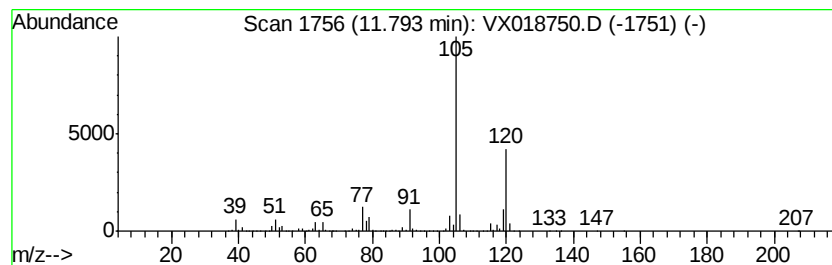
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	17.77 ug/L	2501380	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Mesitylene	120	C9H12	000108-67-8	93
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

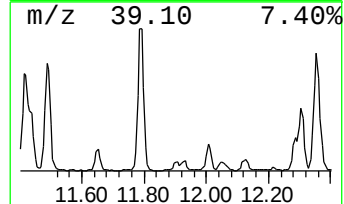
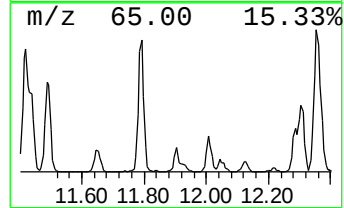
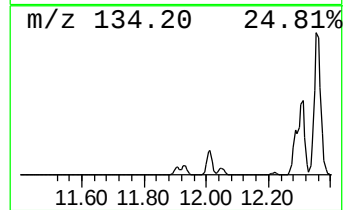
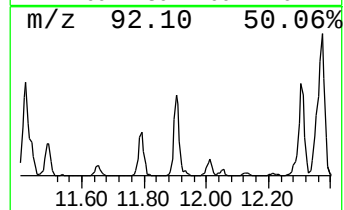
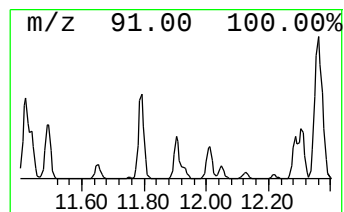
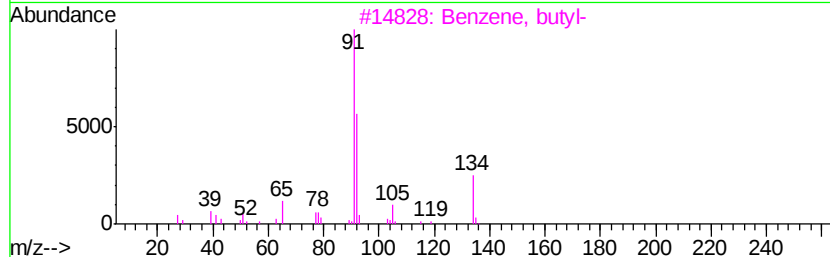
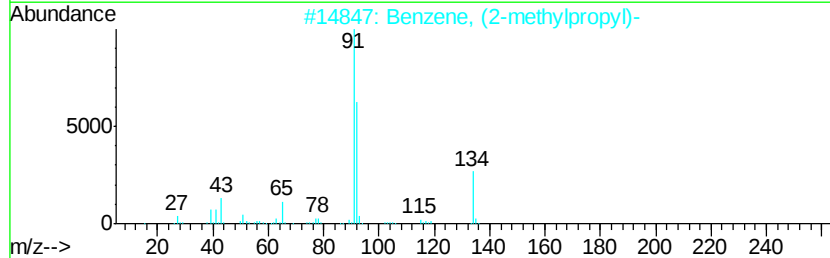
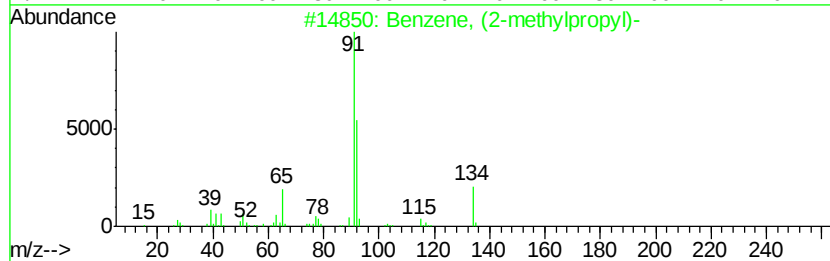
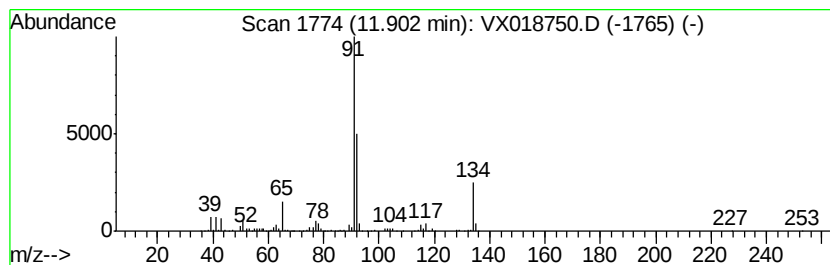
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, (2-methylpropyl)- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	1.12 ug/L	157646	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2		Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	87
3		Benzene, butyl-	134	C10H14	000104-51-8	86
4		Benzene, butyl-	134	C10H14	000104-51-8	80
5		Benzene, butyl-	134	C10H14	000104-51-8	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

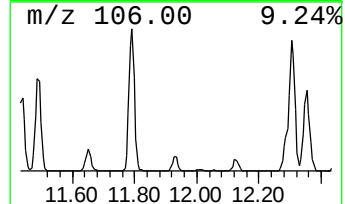
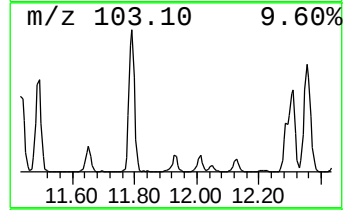
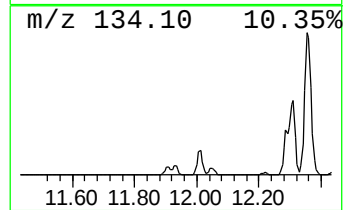
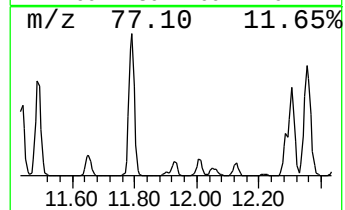
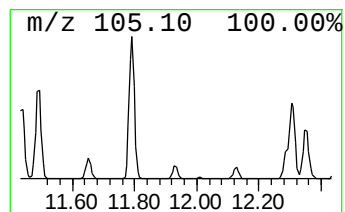
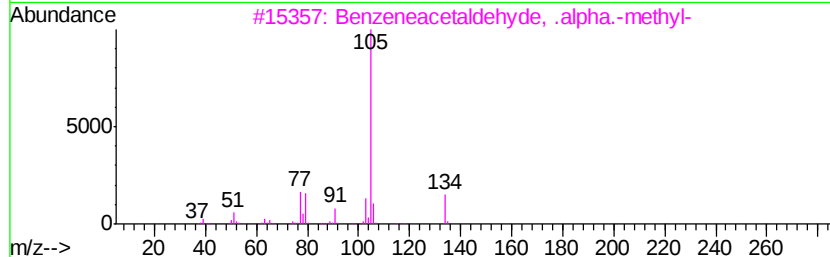
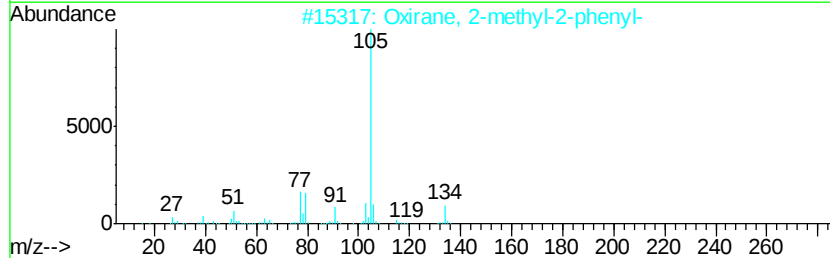
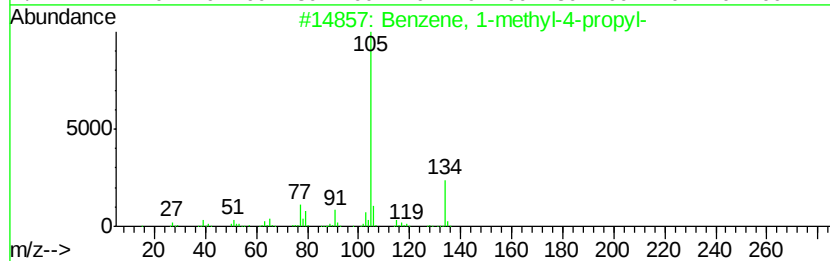
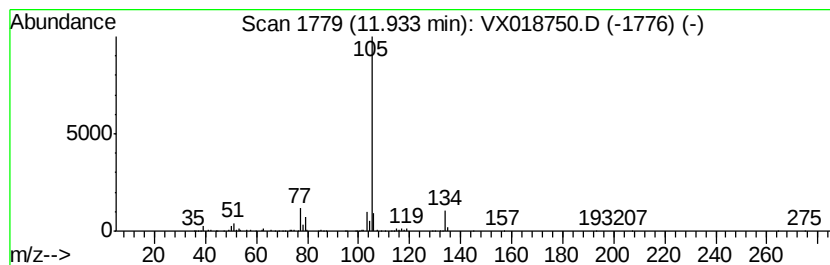
Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Oxirane, 2-methyl-2-phenyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.93	1.32 ug/L	185471	1,4-Dichlorobenzene-d4	12.06	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1		Benzene, 1-methyl-4-propyl-	134 C10H14	001074-55-1	91
2		Oxirane, 2-methyl-2-phenyl-	134 C9H10O	002085-88-3	90
3		Benzeneacetaldehyde, .alpha.-met...	134 C9H10O	000093-53-8	90
4		Benzene, (1-methylpropyl)-	134 C10H14	000135-98-8	86
5		Benzeneacetaldehyde, 4-methyl-	134 C9H10O	000104-09-6	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

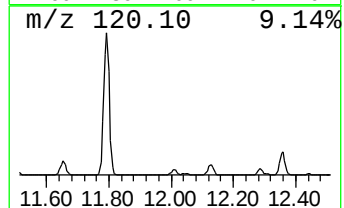
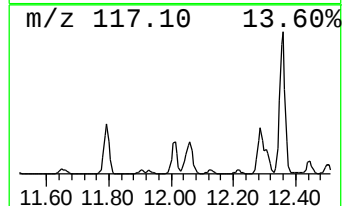
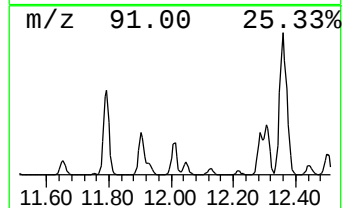
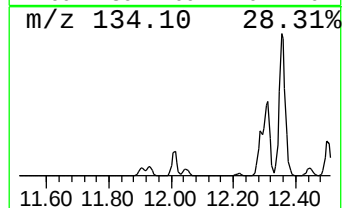
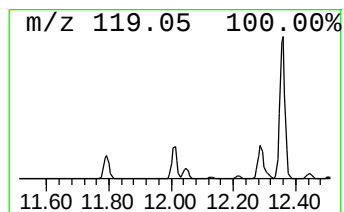
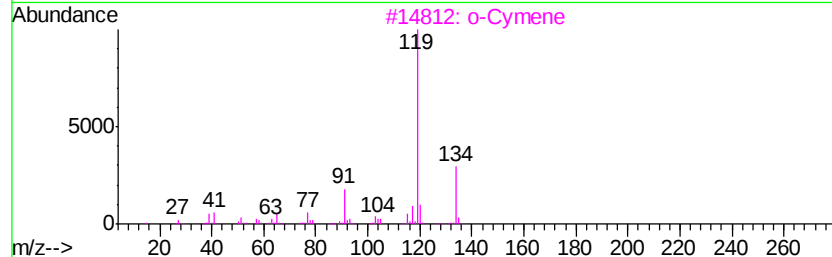
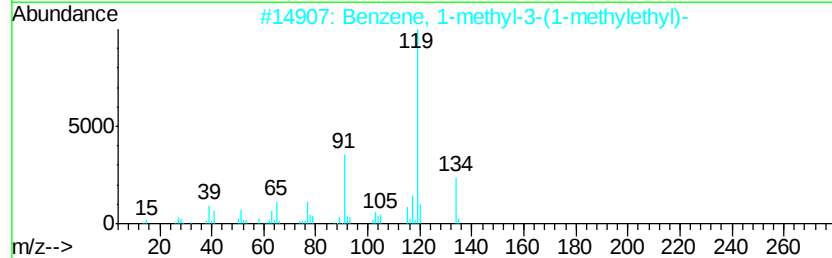
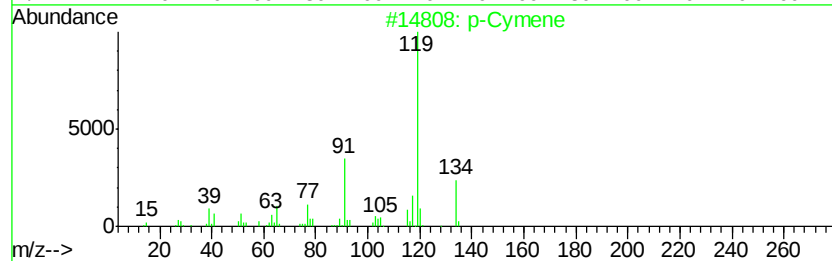
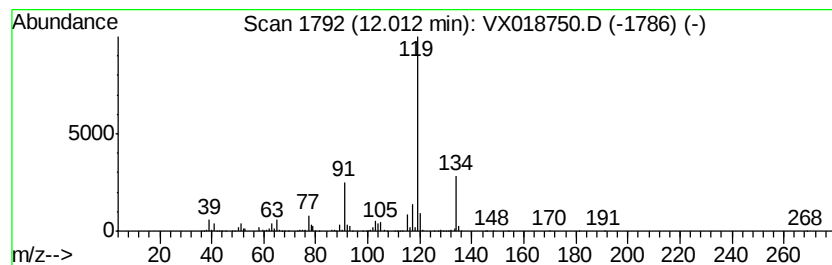
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Benzene, 1-methyl-3-(1-meth... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	3.02 ug/L	424472	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	95
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3		o-Cymene	134	C10H14	000527-84-4	94
4		p-Cymene	134	C10H14	000099-87-6	94
5		o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

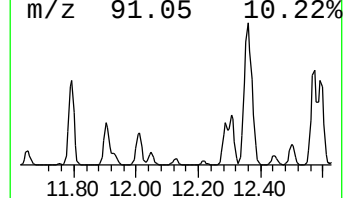
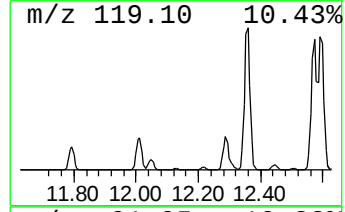
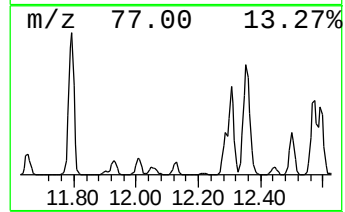
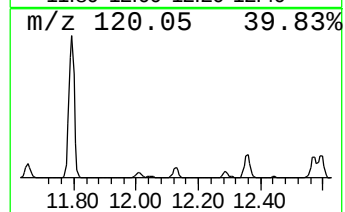
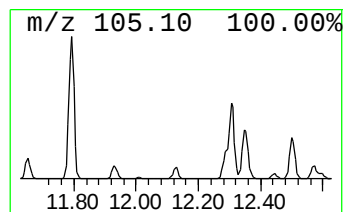
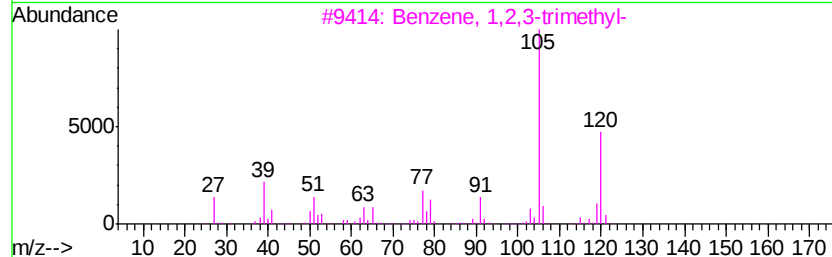
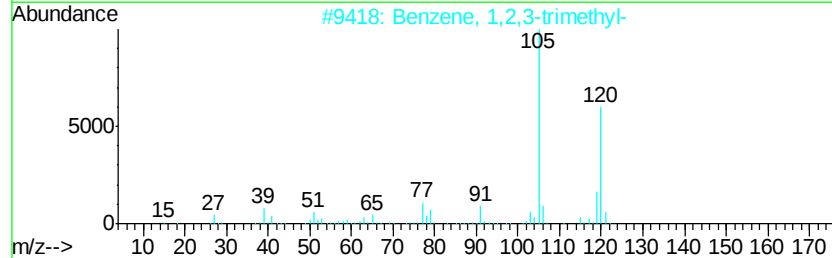
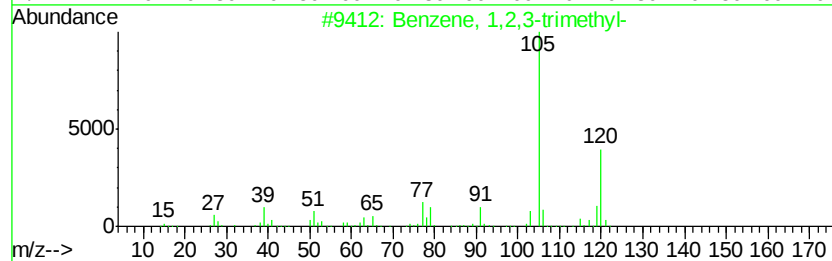
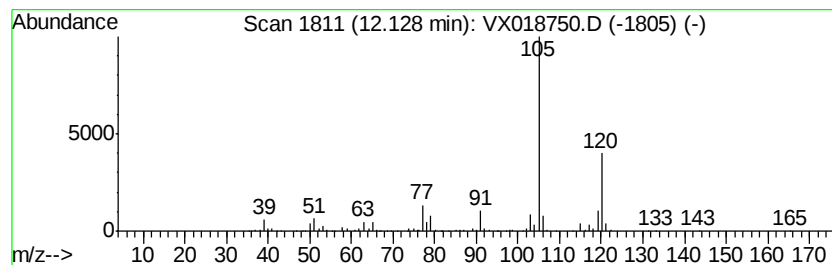
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Benzene, 1,2,4-trimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	1.42 ug/L	199354	1,4-Dichlorobenzene-d4	12.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

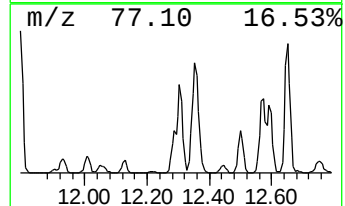
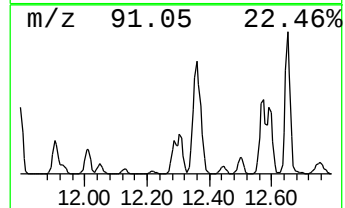
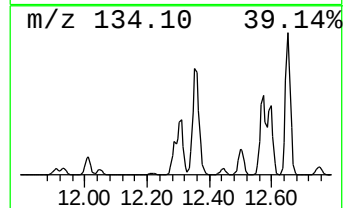
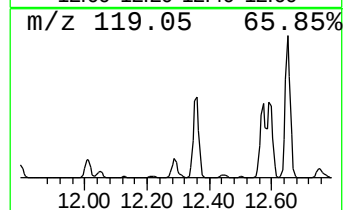
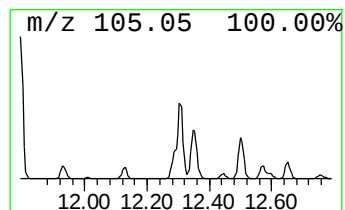
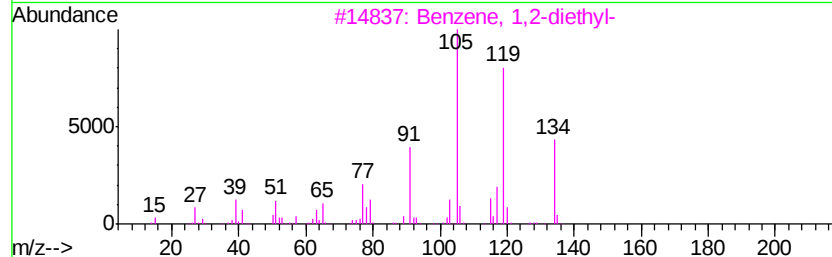
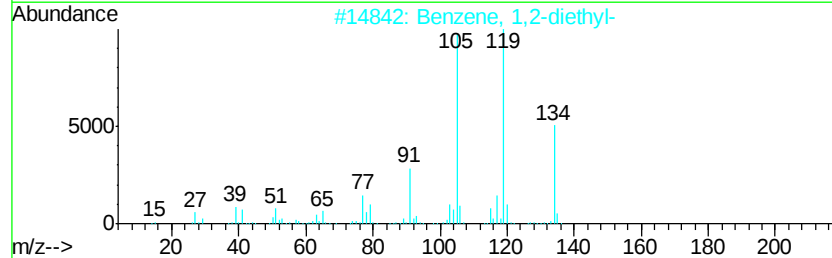
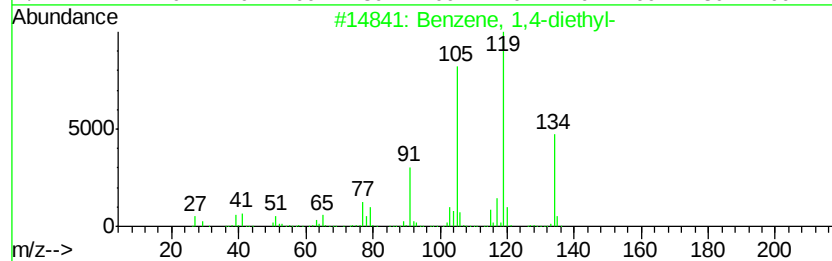
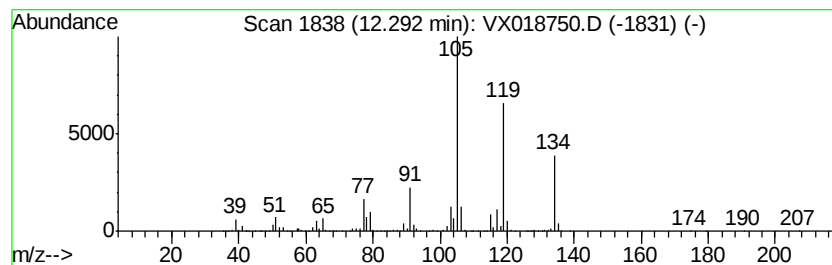
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Benzene, 1,4-diethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	5.17 ug/L	727662	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
4		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
5		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

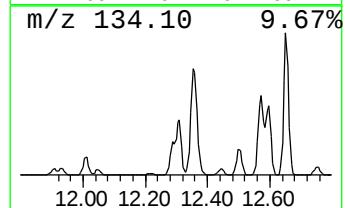
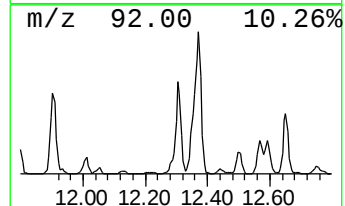
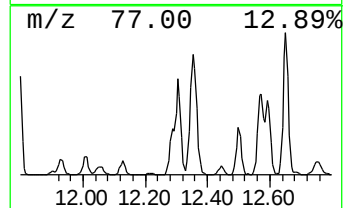
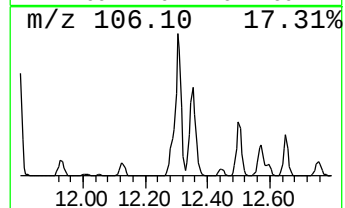
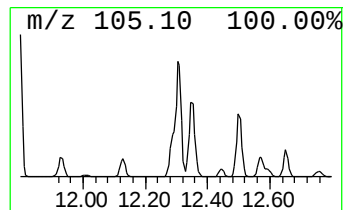
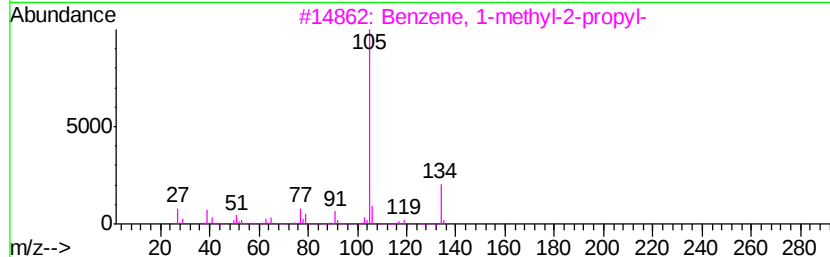
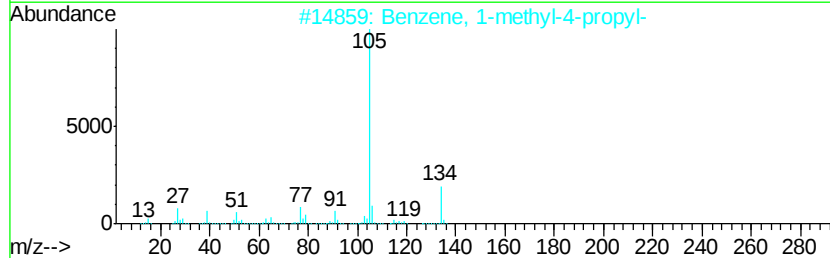
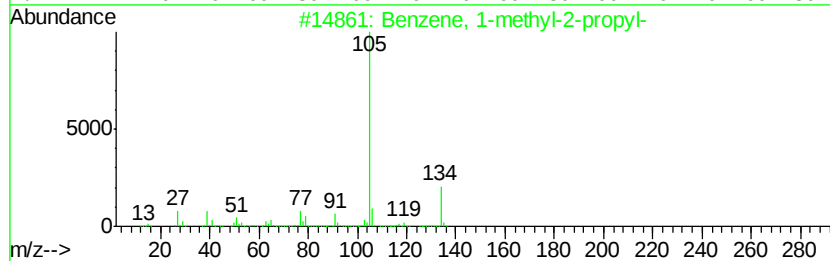
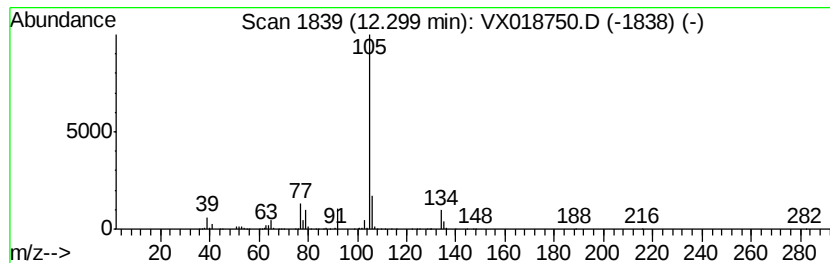
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzene, 1-methyl-2-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	8.48 ug/L	1194350	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	87
2		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	86
3		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	80
4		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	78
5		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

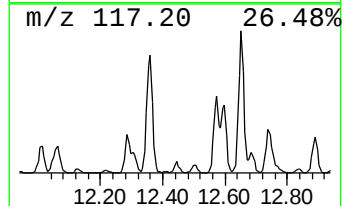
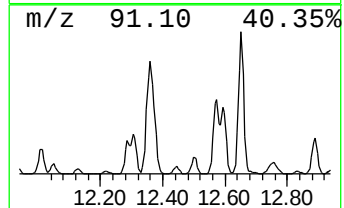
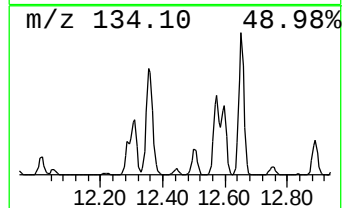
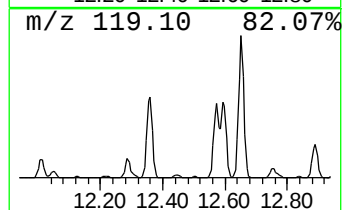
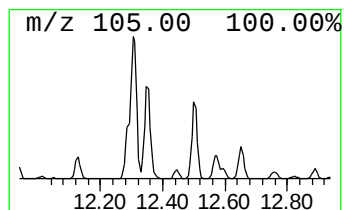
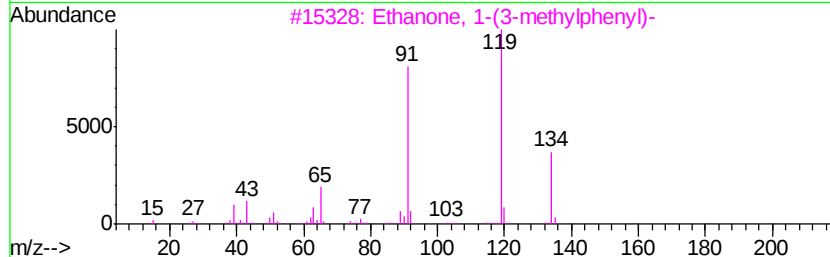
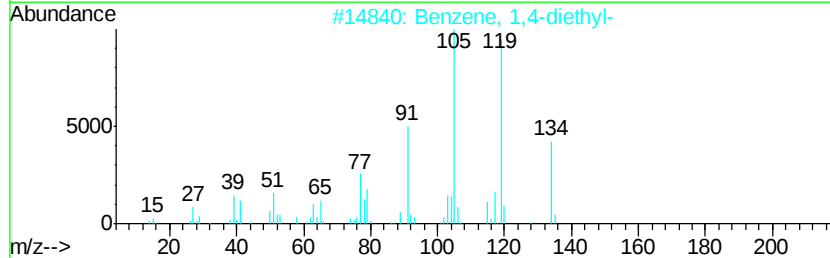
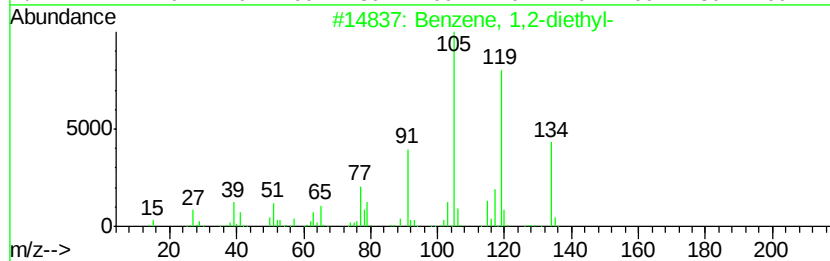
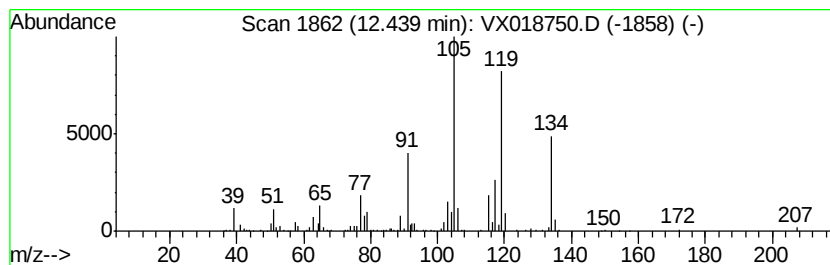
Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Benzene, 1,2-diethyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.44	1.22 ug/L	171526	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 96
2		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 95
3		Ethanone, 1-(3-methylphenyl)-	134 C9H10O	000585-74-0 91
4		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 90
5		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

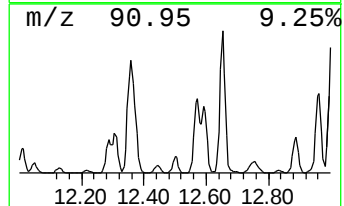
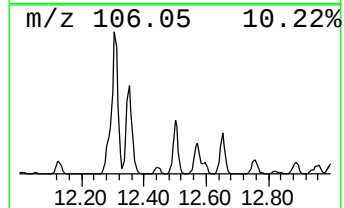
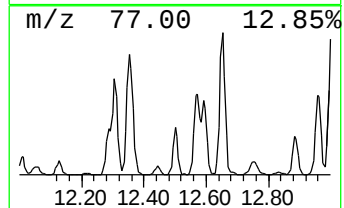
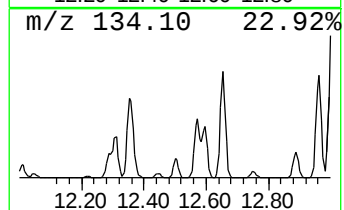
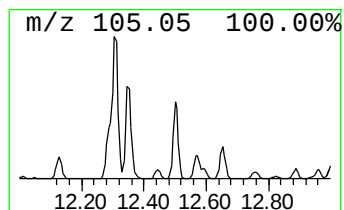
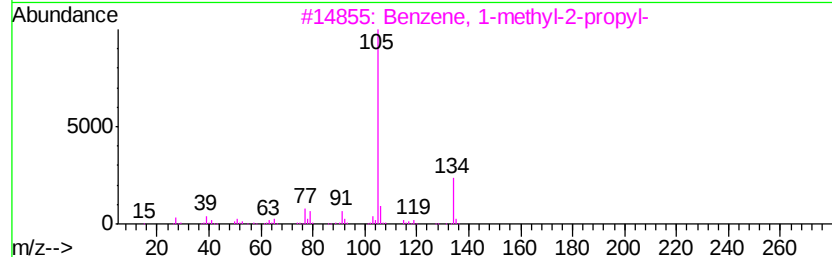
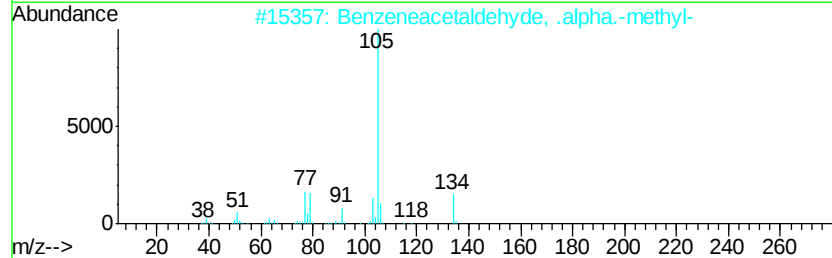
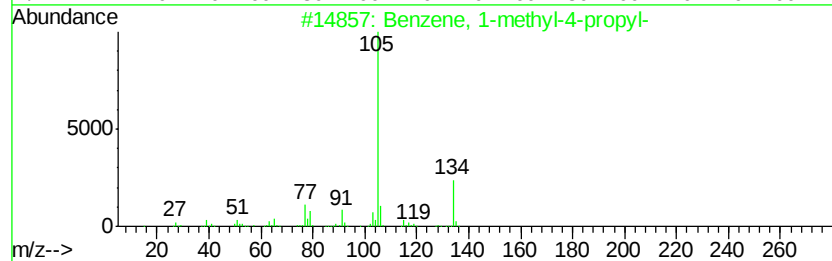
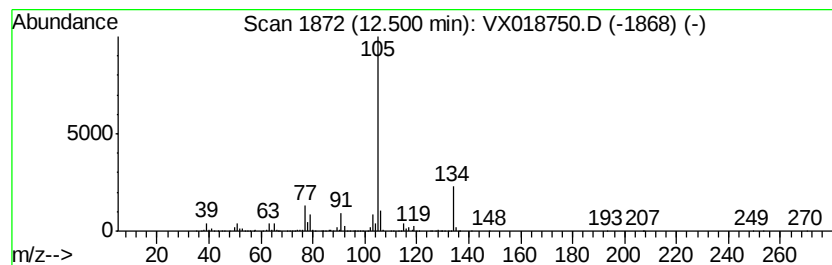
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 Benzeneacetaldehyde, .alpha... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.50	4.34 ug/L	610350	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	91
3		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

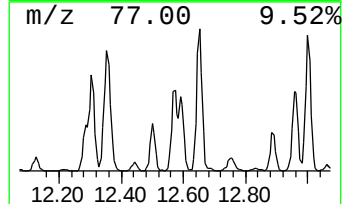
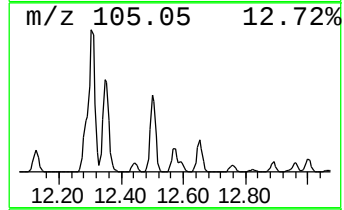
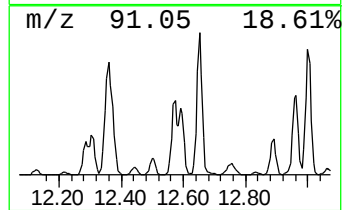
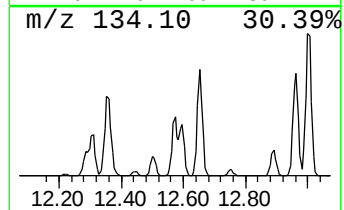
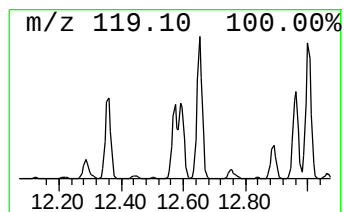
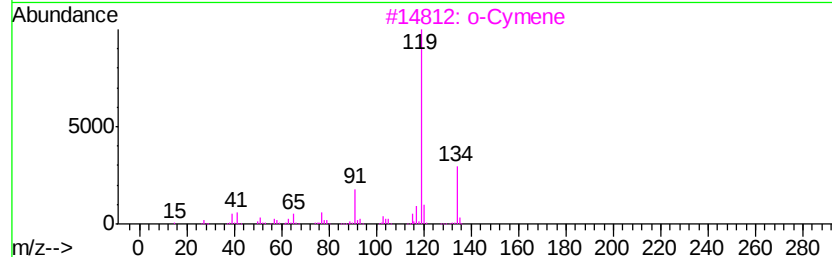
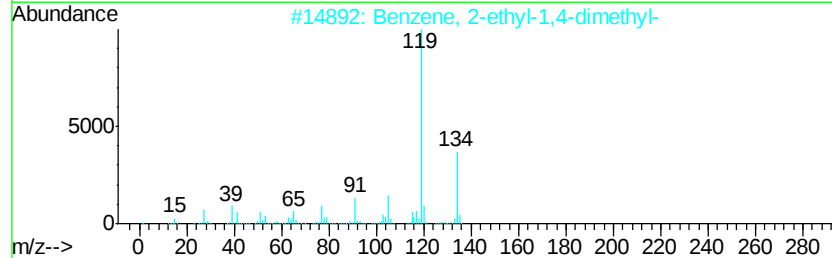
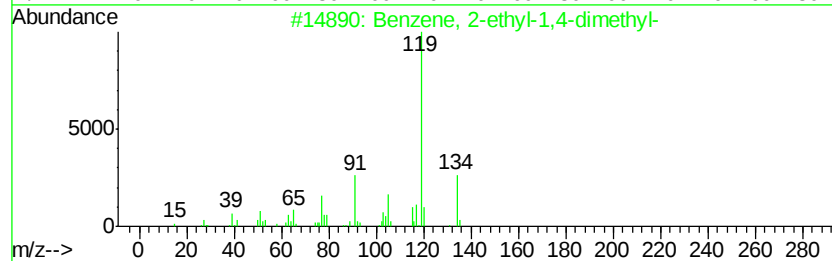
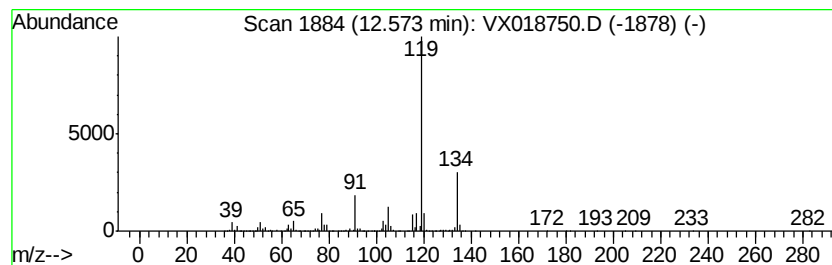
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	13.70 ug/L	1928740	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

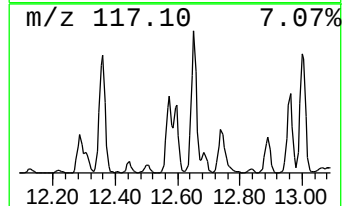
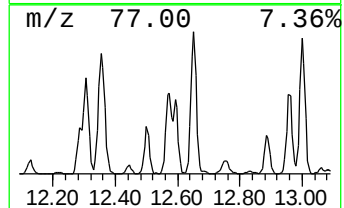
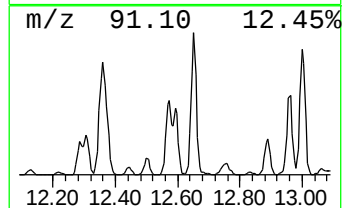
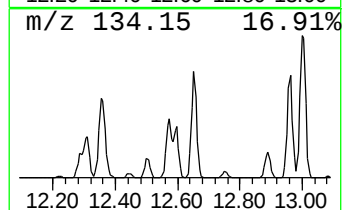
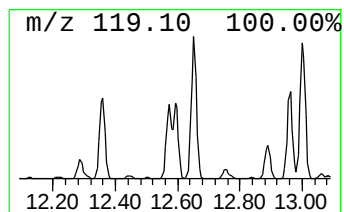
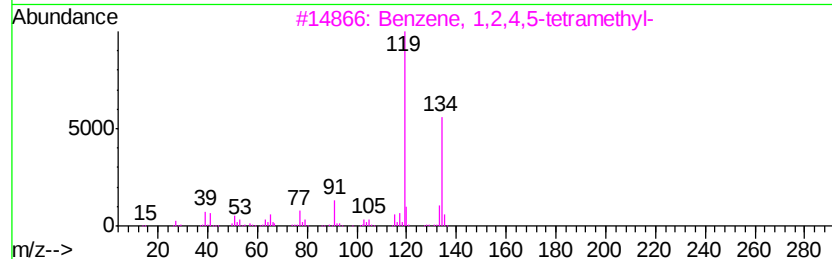
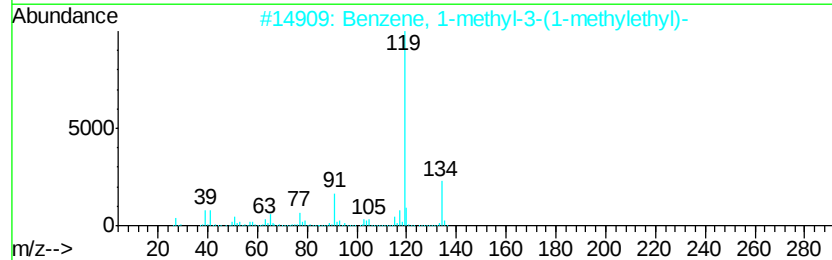
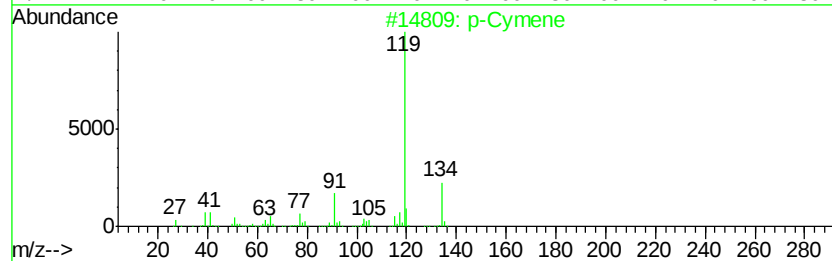
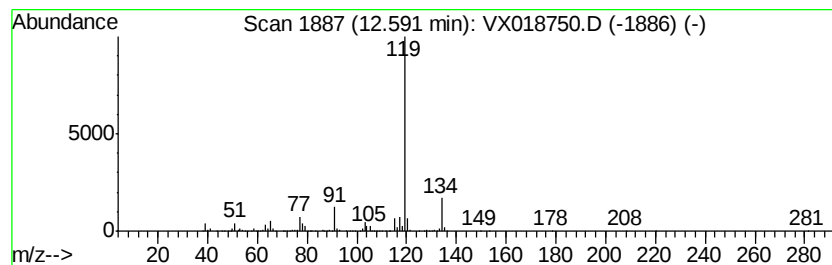
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	8.50 ug/L	1196200	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	95
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
5		o-Cymene	134	C10H14	000527-84-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

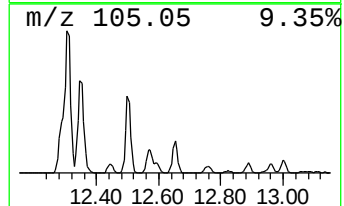
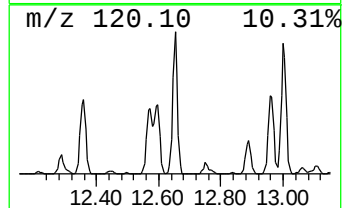
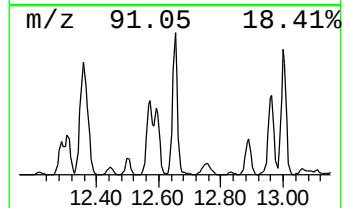
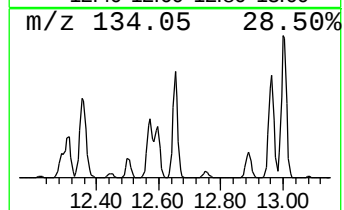
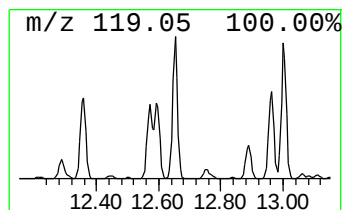
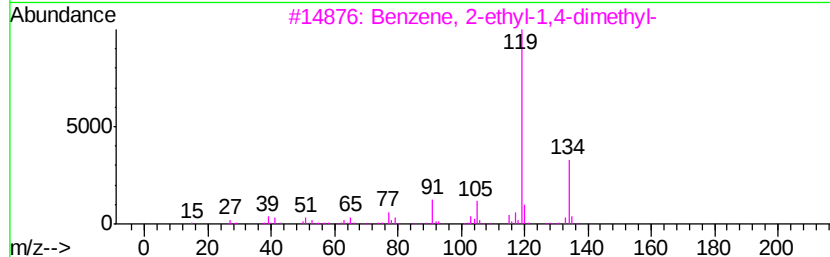
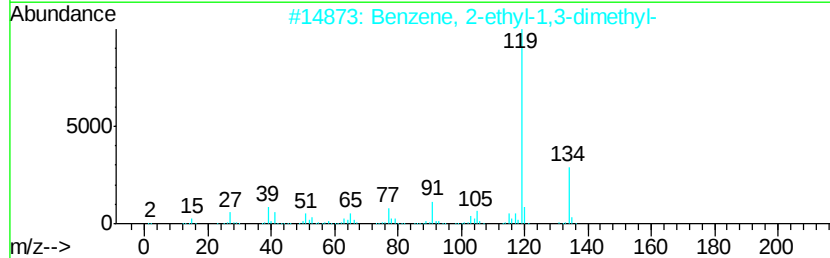
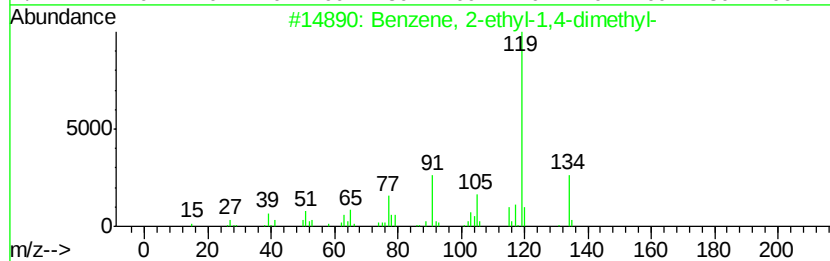
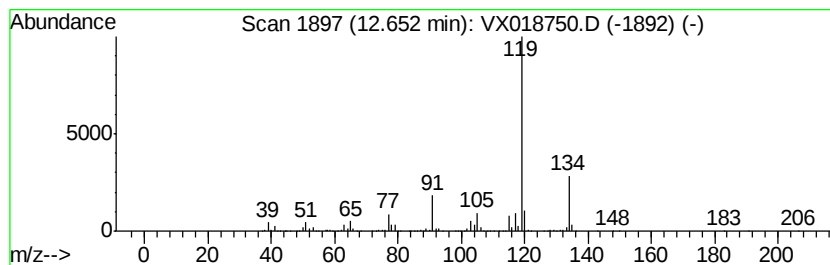
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	20.75 ug/L	2921590	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

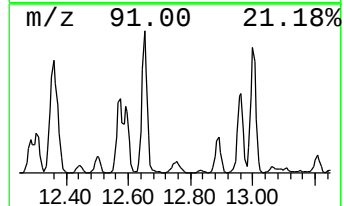
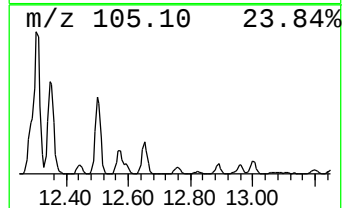
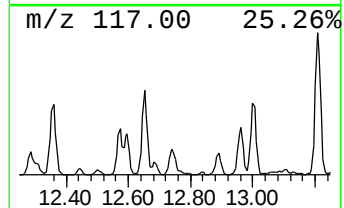
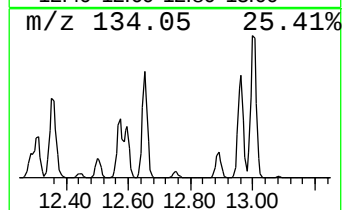
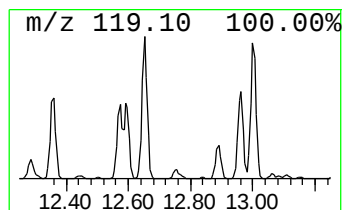
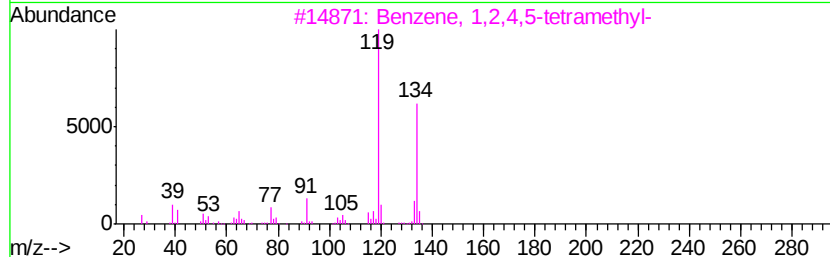
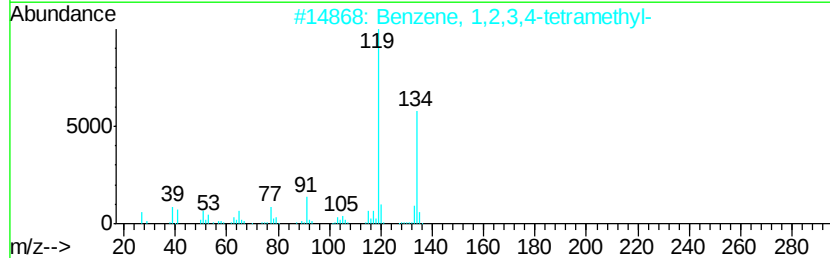
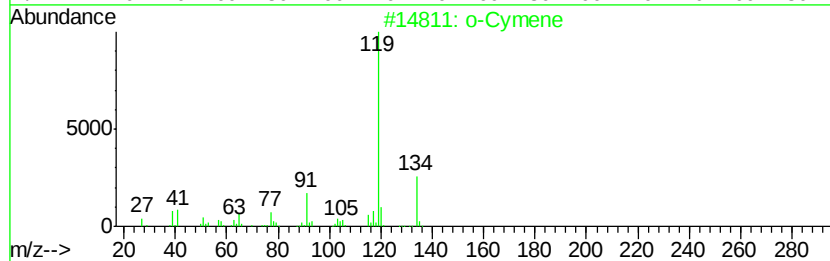
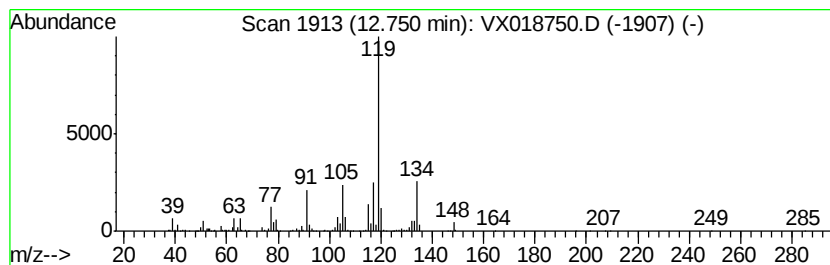
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 o-Cymene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	2.89 ug/L	406767	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	92
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	76
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	76
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	76
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

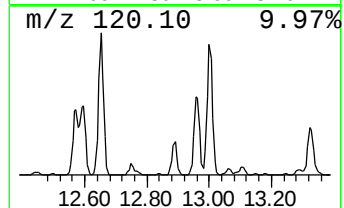
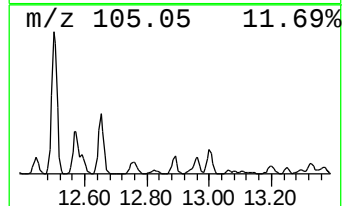
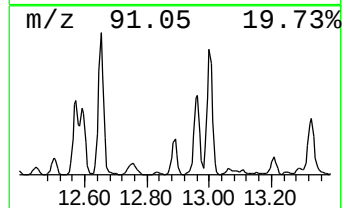
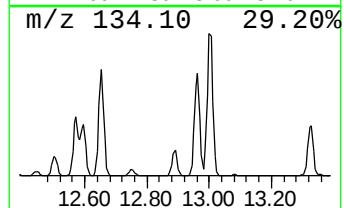
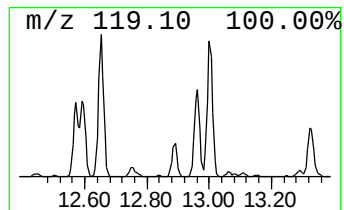
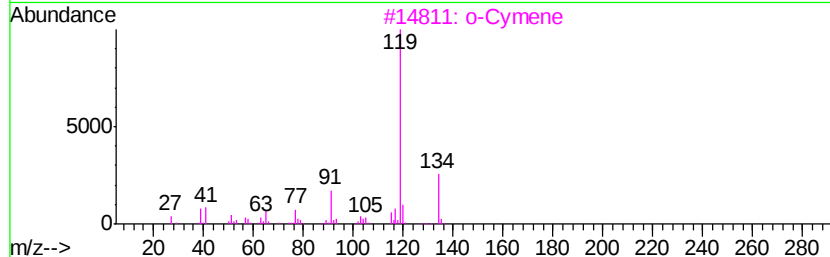
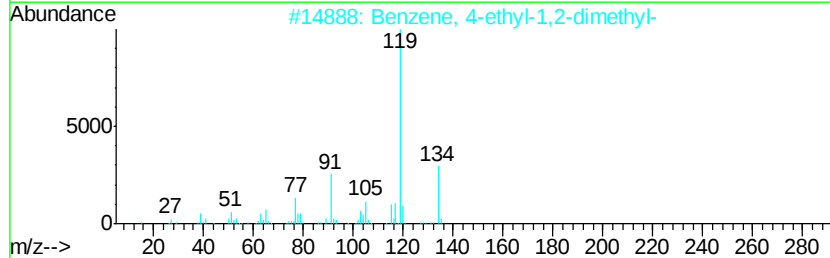
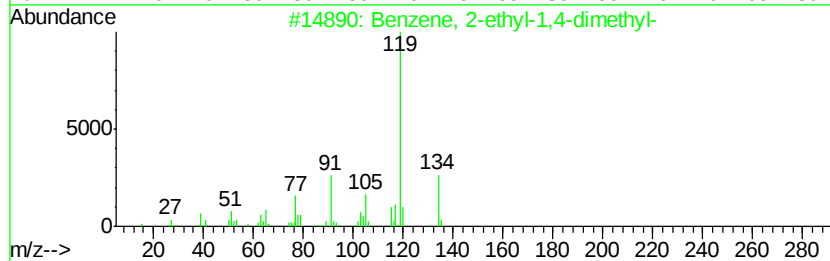
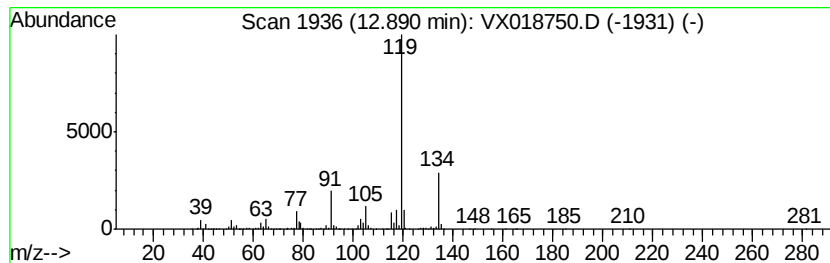
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	5.46 ug/L	768565	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

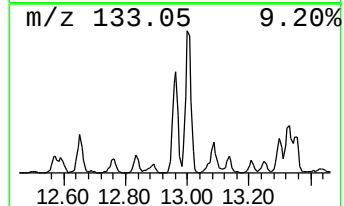
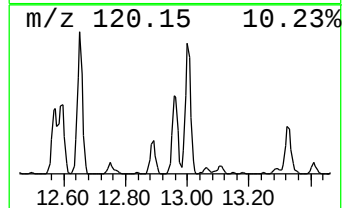
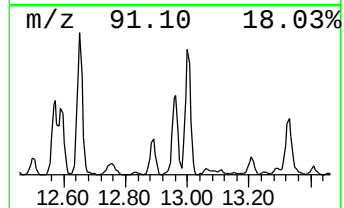
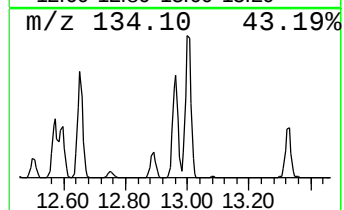
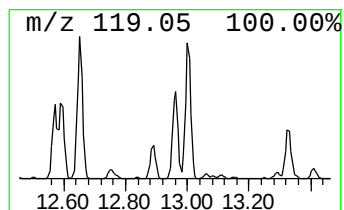
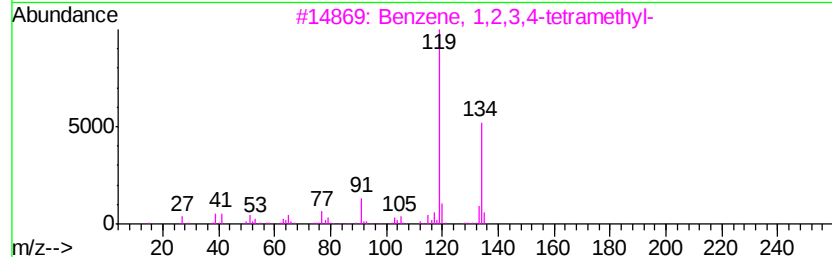
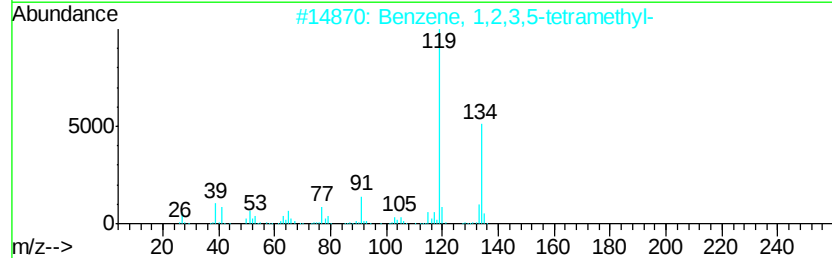
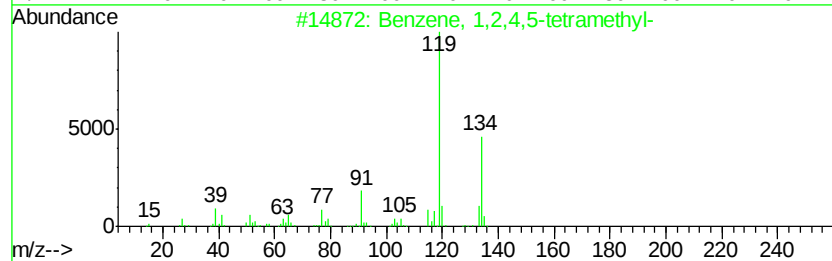
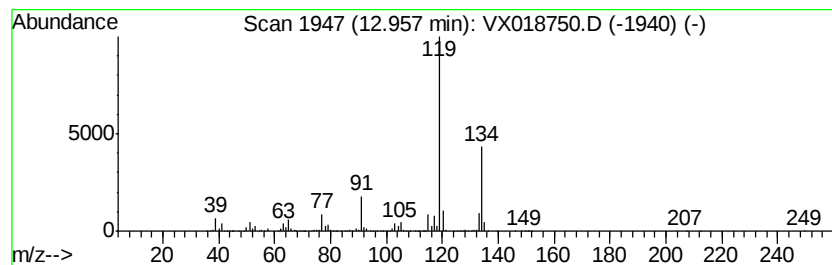
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	14.36 ug/L	2021330	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

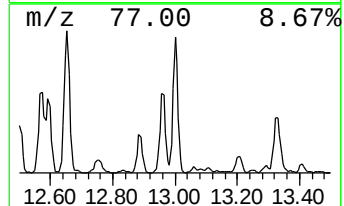
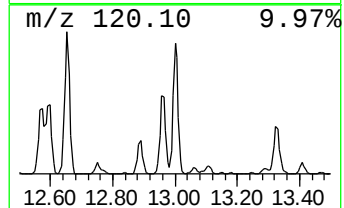
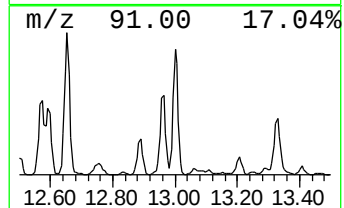
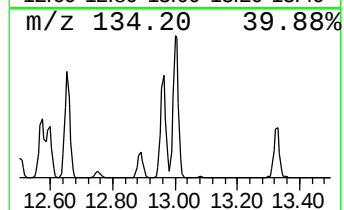
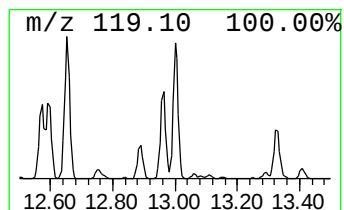
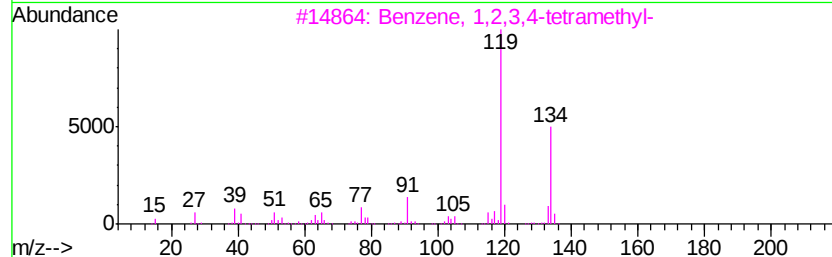
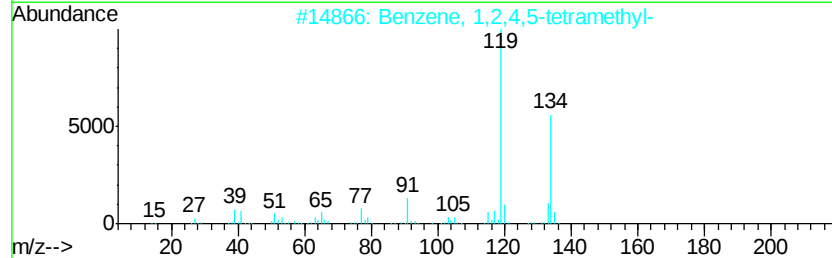
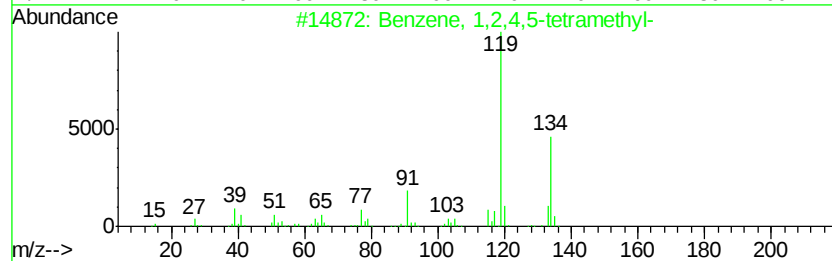
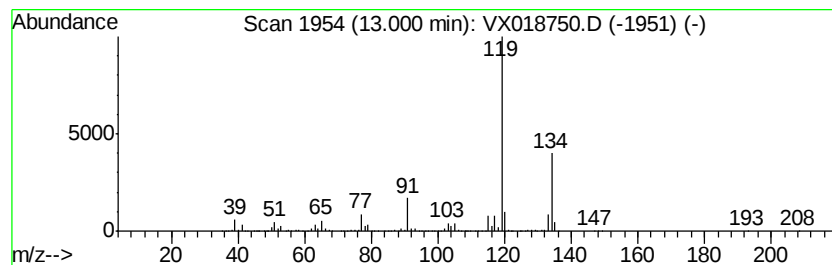
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	21.48 ug/L	3023810	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

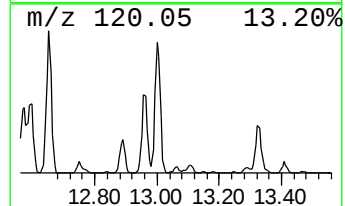
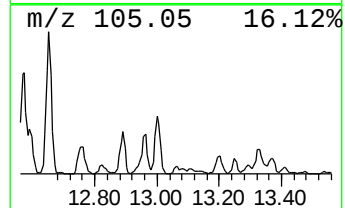
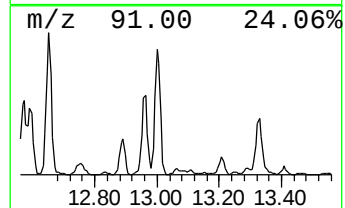
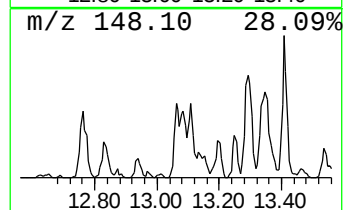
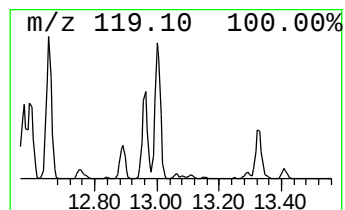
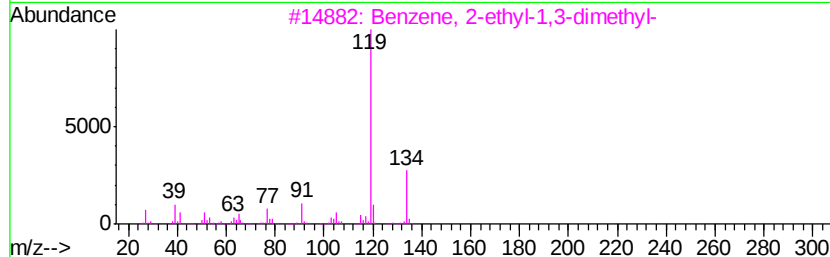
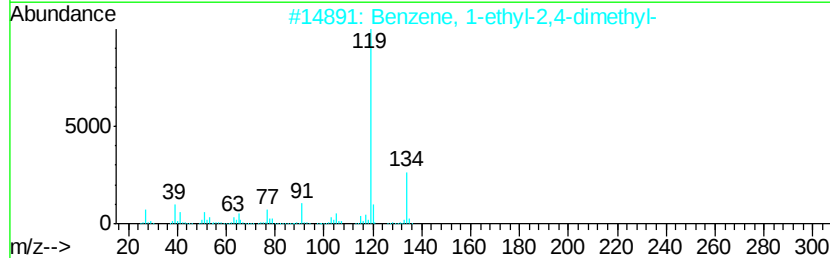
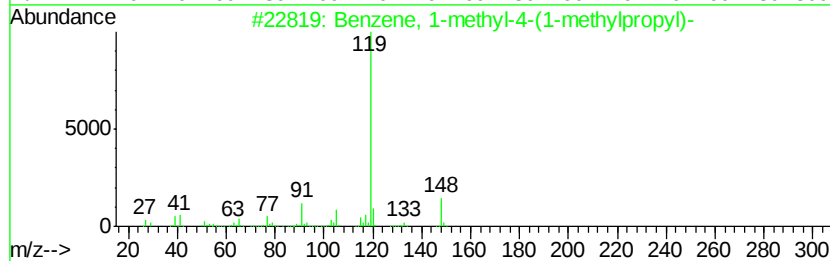
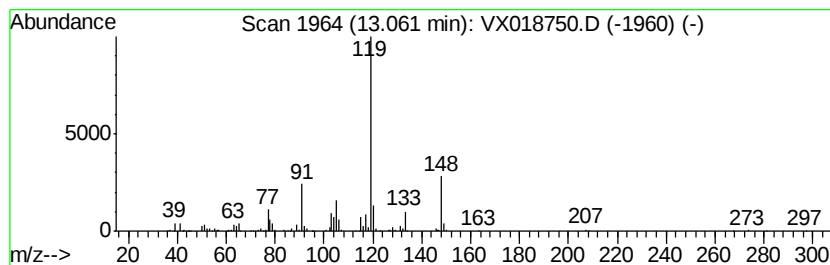
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 Benzene, 1-methyl-4-(1-meth... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.06	0.76 ug/L	107056	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	87
2		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	68
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	68
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	64
5		2,4-Dimethylphenethyl alcohol	150	C10H14O	006597-59-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

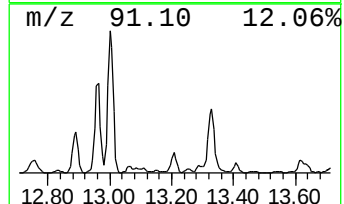
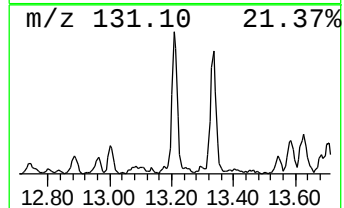
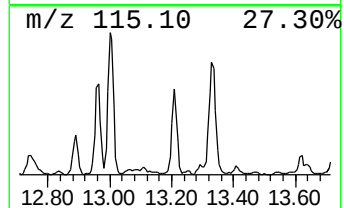
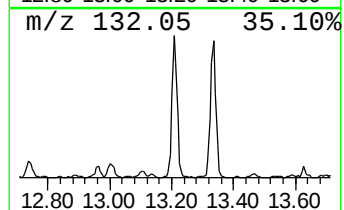
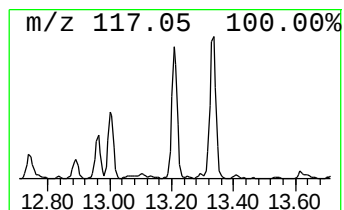
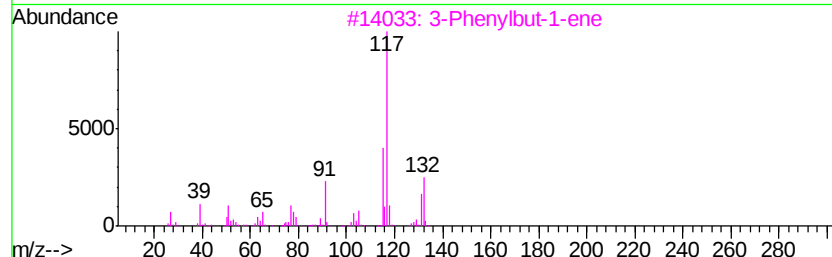
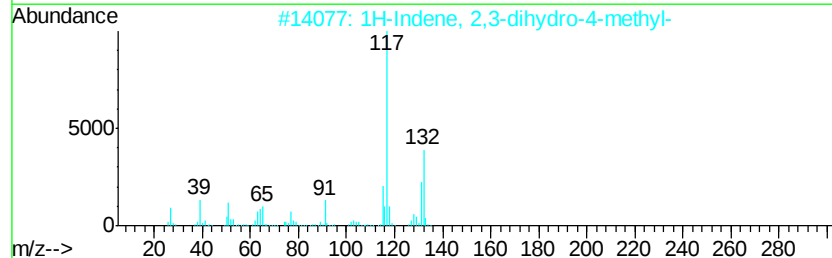
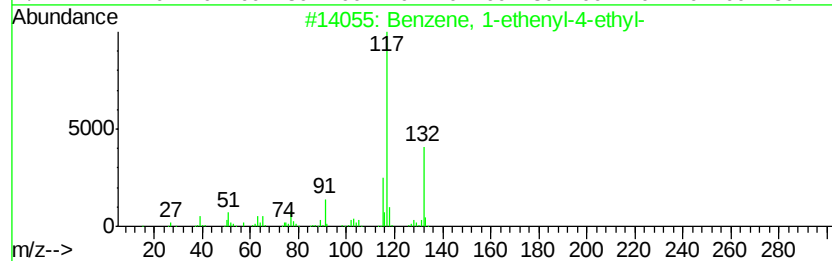
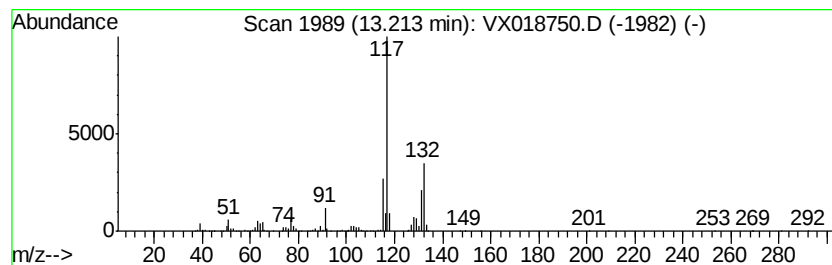
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	4.29 ug/L	604576	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

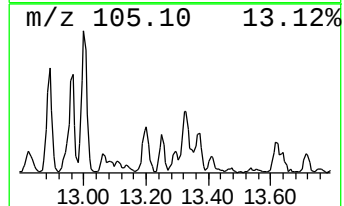
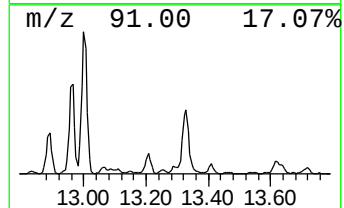
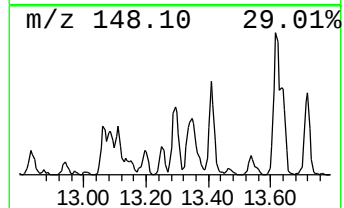
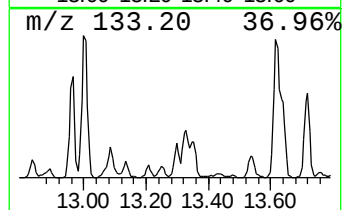
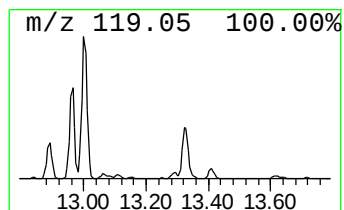
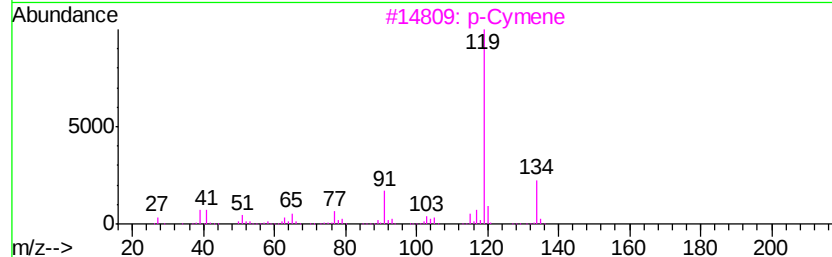
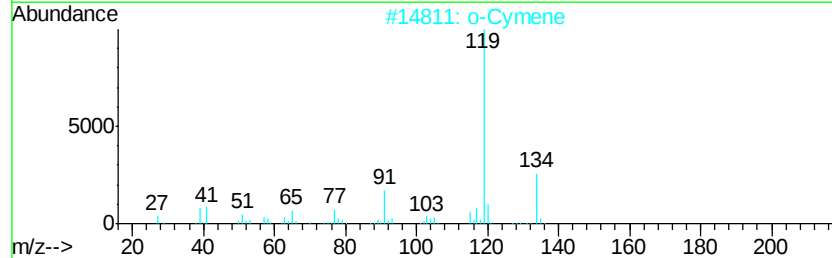
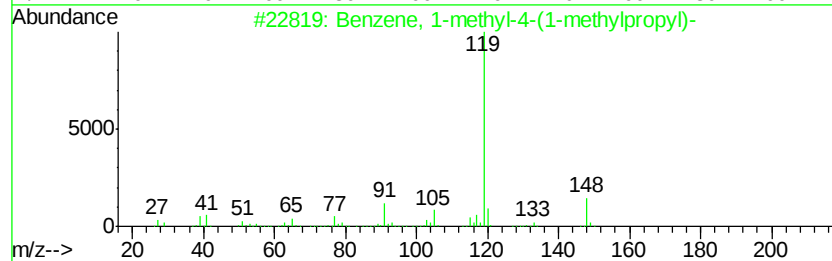
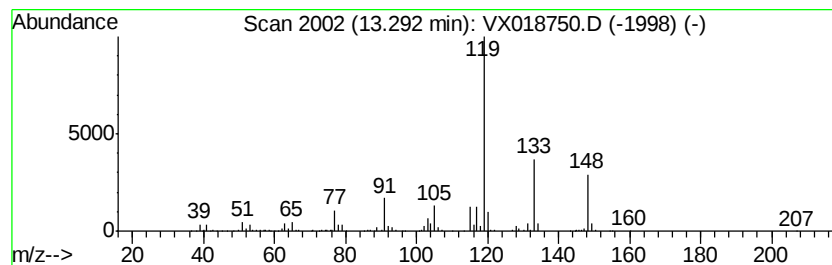
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 unknown-01 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	1.29 ug/L	181541	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	76
2		o-Cymene	134	C10H14	000527-84-4	64
3		p-Cymene	134	C10H14	000099-87-6	64
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

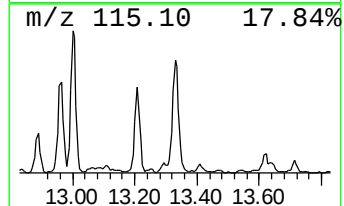
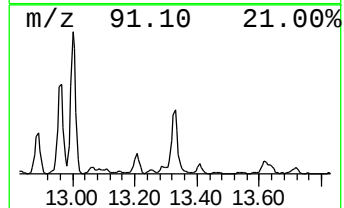
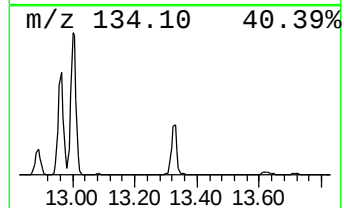
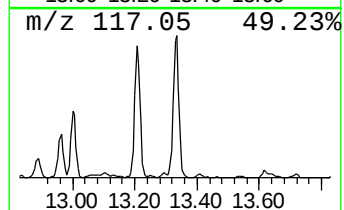
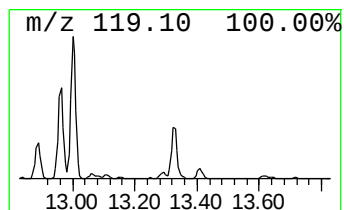
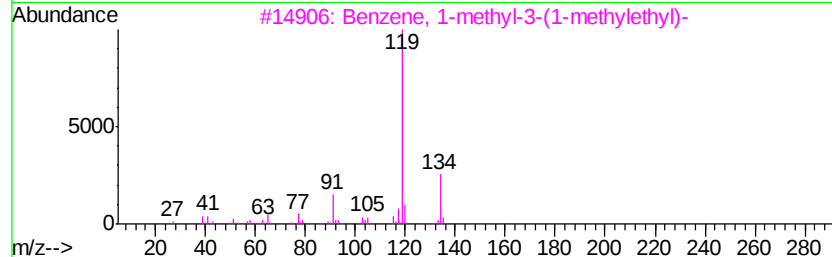
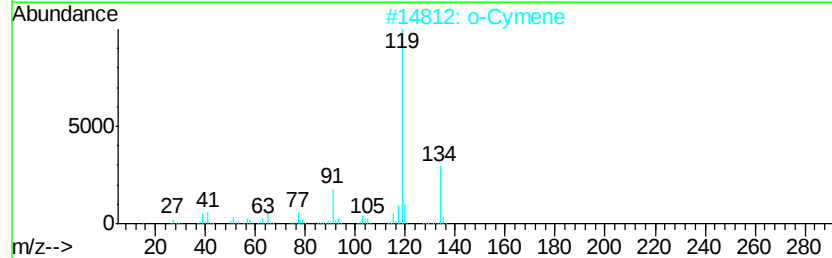
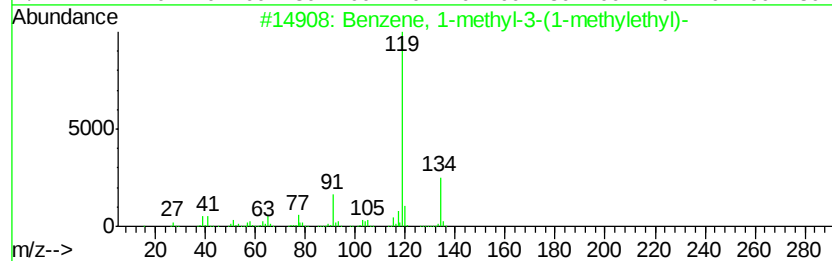
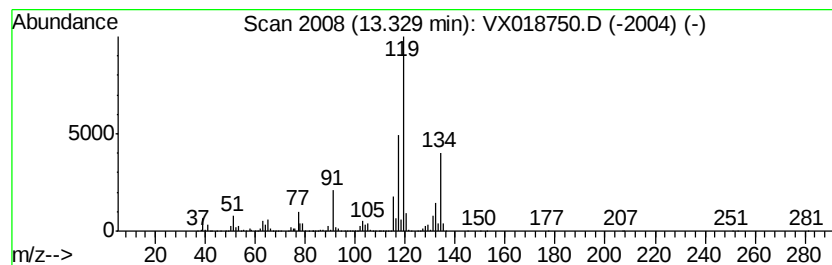
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	12.89 ug/L	1814540	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	70
2		o-Cymene	134	C10H14	000527-84-4	70
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	70
4		o-Cymene	134	C10H14	000527-84-4	70
5		p-Cymene	134	C10H14	000099-87-6	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

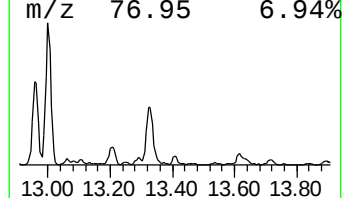
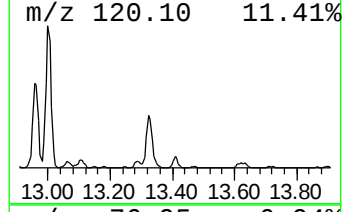
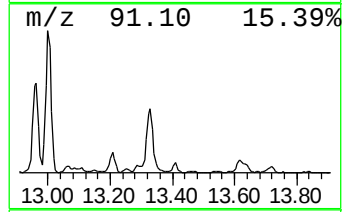
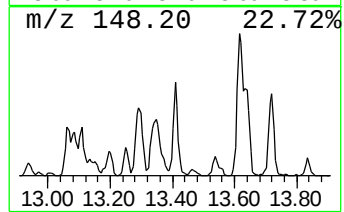
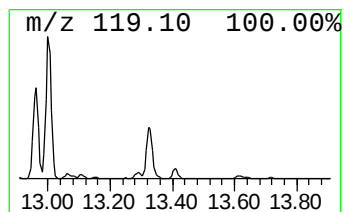
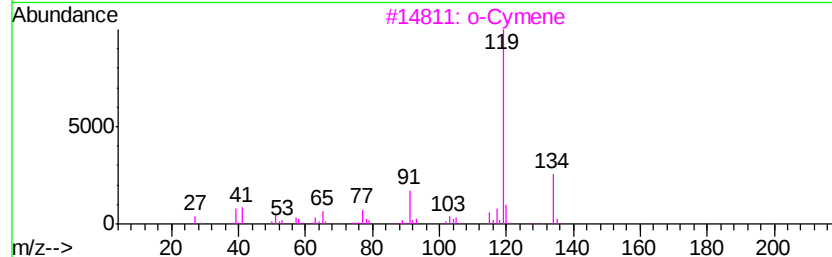
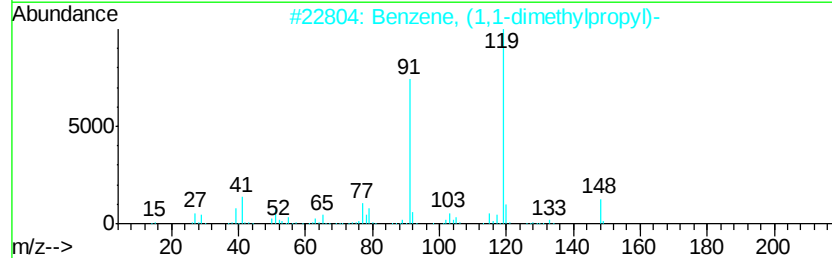
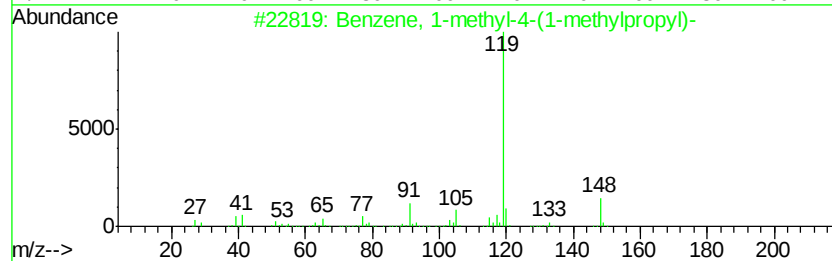
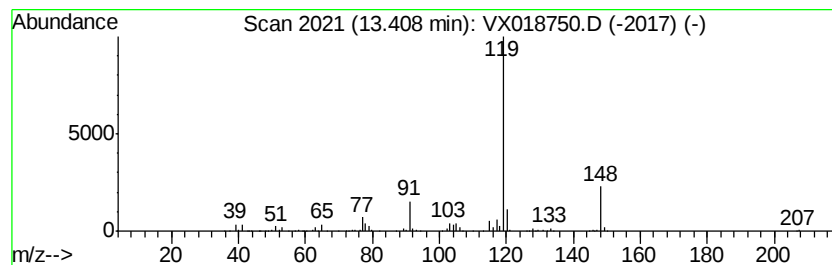
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 Benzene, (1,1-dimethylpropyl)- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	1.24 ug/L	174212	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	91
2		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	86
3		o-Cymene	134	C10H14	000527-84-4	72
4		p-Cymene	134	C10H14	000099-87-6	72
5		o-Cymene	134	C10H14	000527-84-4	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

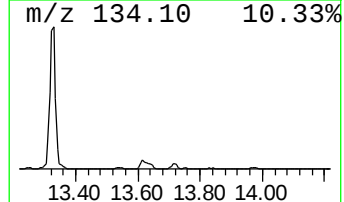
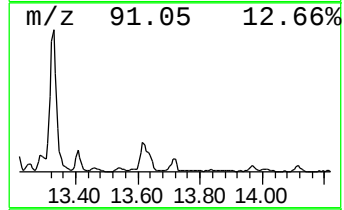
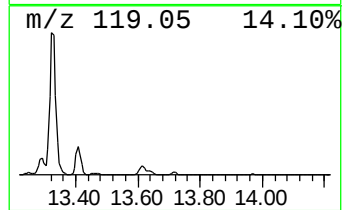
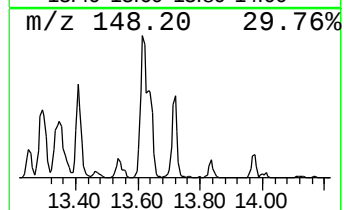
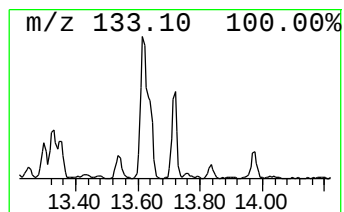
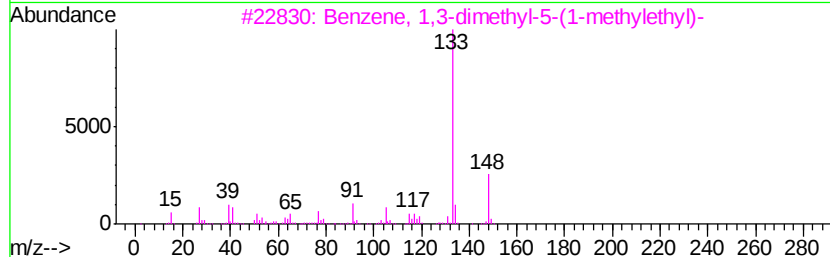
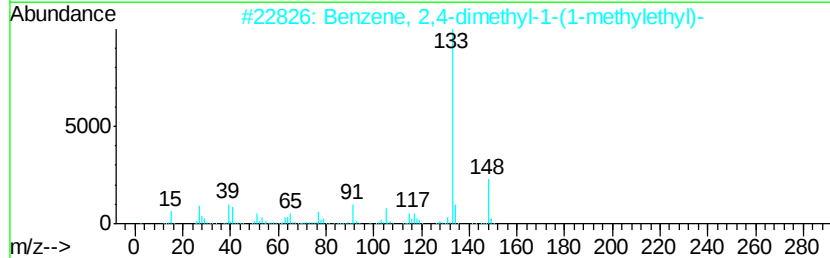
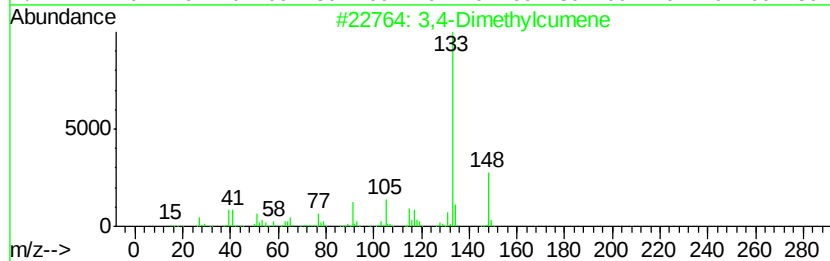
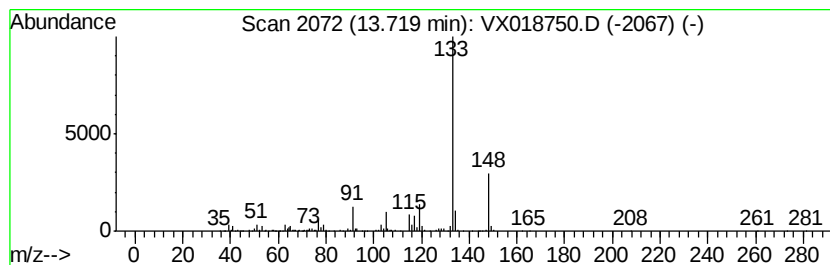
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 3,4-Dimethylcumene Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	1.20 ug/L	168860	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,4-Dimethylcumene	148	C11H16	1000370-34-1	90
2		Benzene, 2,4-dimethyl-1-(1-methyl-...	148	C11H16	004706-89-2	90
3		Benzene, 1,3-dimethyl-5-(1-methyl-...	148	C11H16	004706-90-5	90
4		Benzene, 1,4-dimethyl-2-(1-methyl-...	148	C11H16	004132-72-3	87
5		Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	83



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

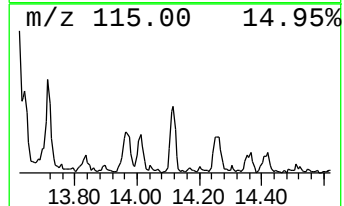
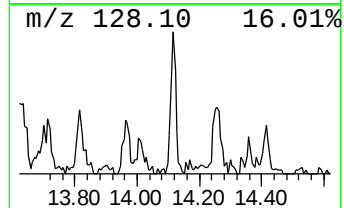
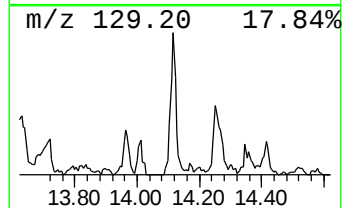
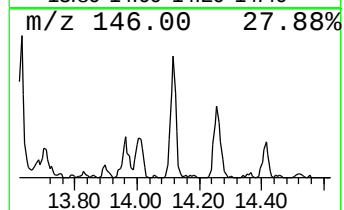
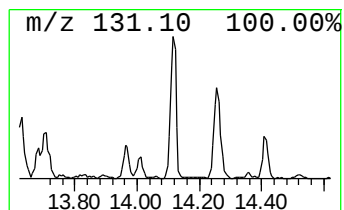
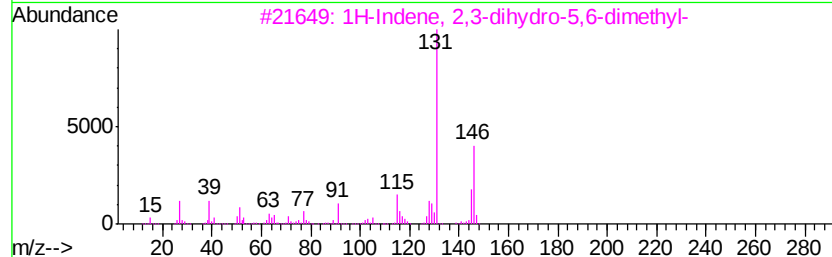
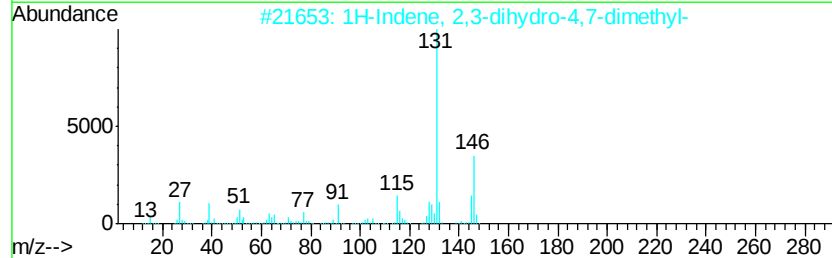
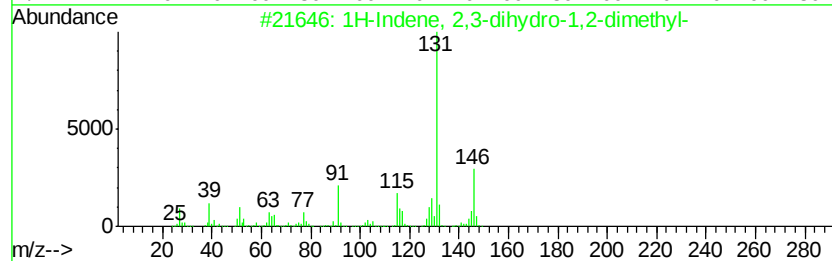
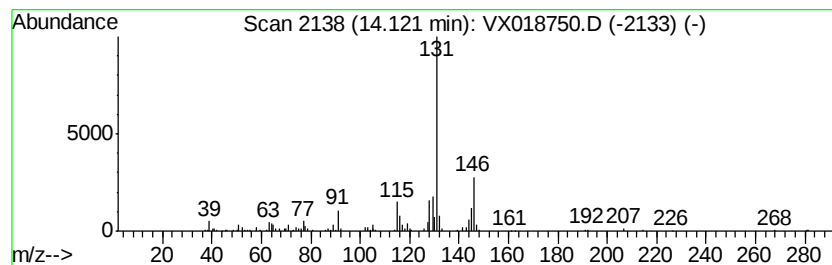
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.12	0.70 ug/L	98188	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
3		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	81
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3W0DL2

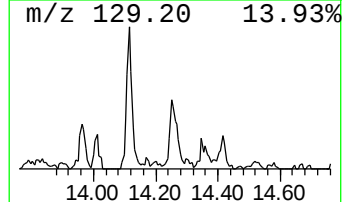
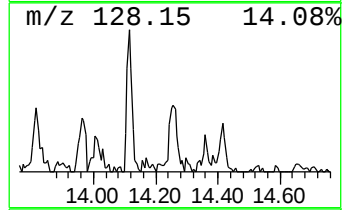
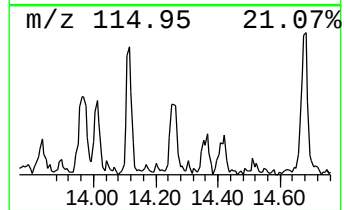
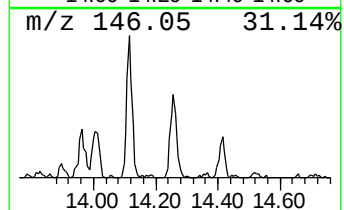
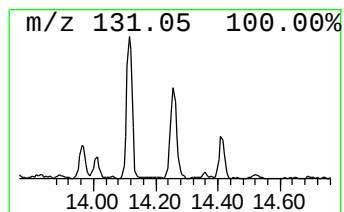
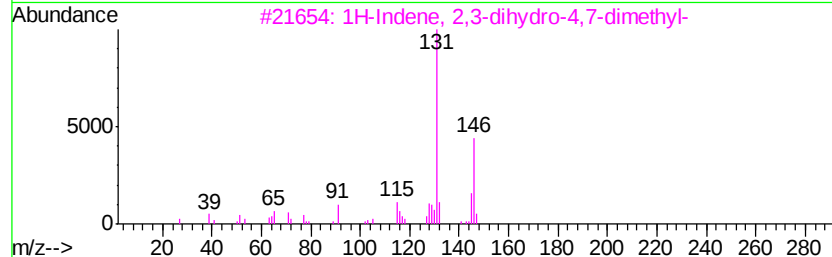
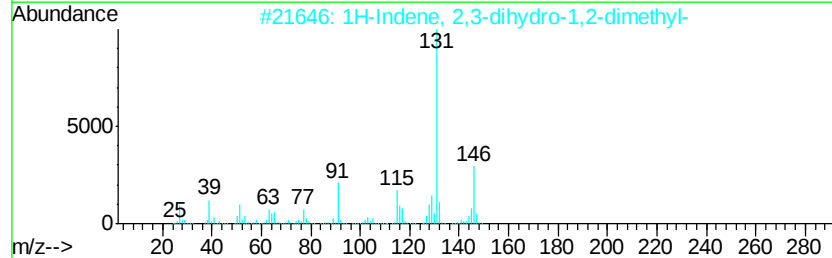
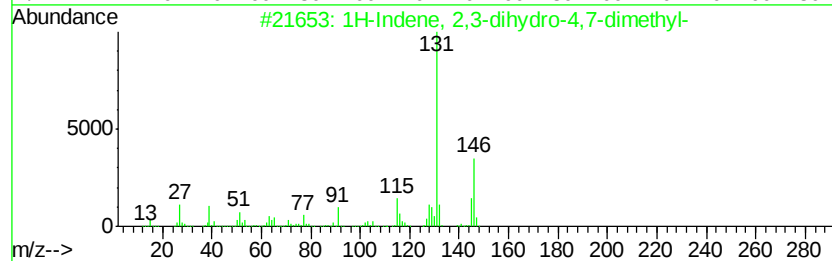
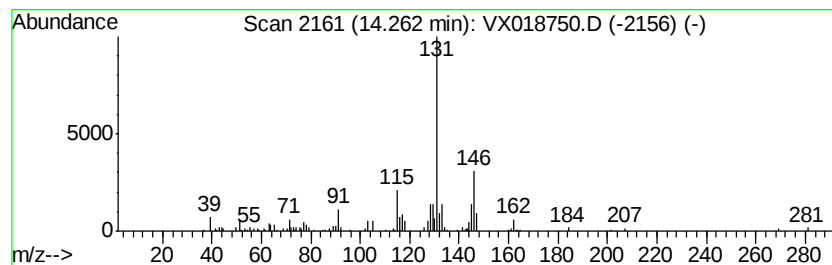
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	0.55 ug/L	76843	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	93
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4		Benzene, (3-methyl-2-butenvl)-	146	C11H14	004489-84-3	91
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100220\
Data File : VX018750.D
Acq On : 02 Oct 2020 16:16
Operator : JC/SP
Sample : L4171-13DL2 50X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3W0DL2

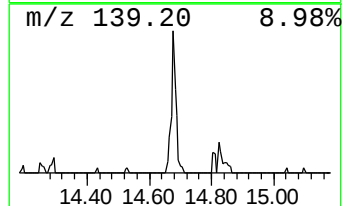
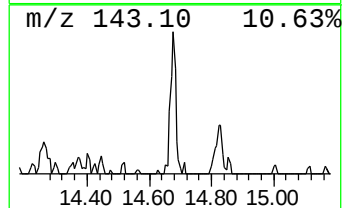
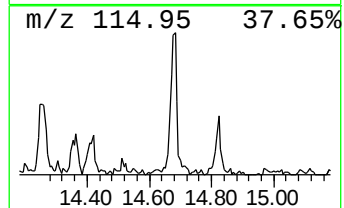
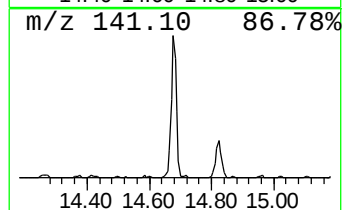
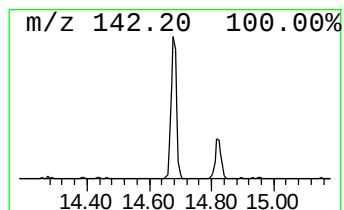
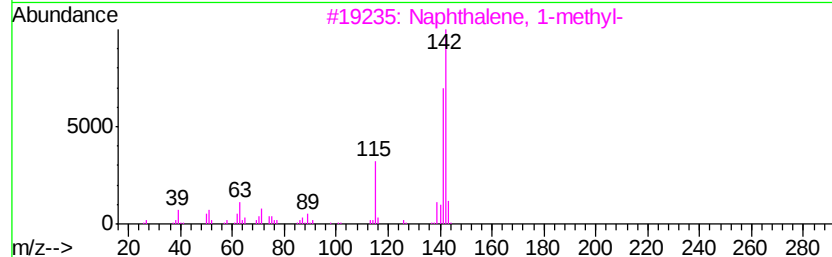
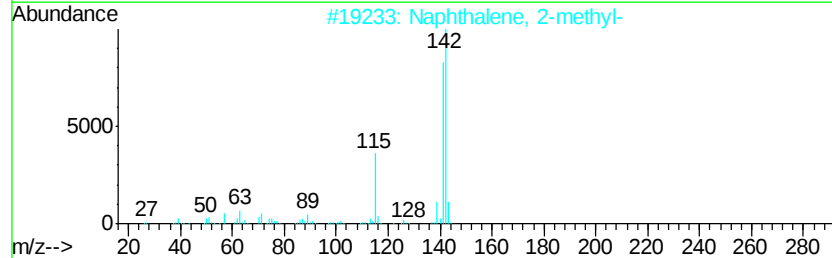
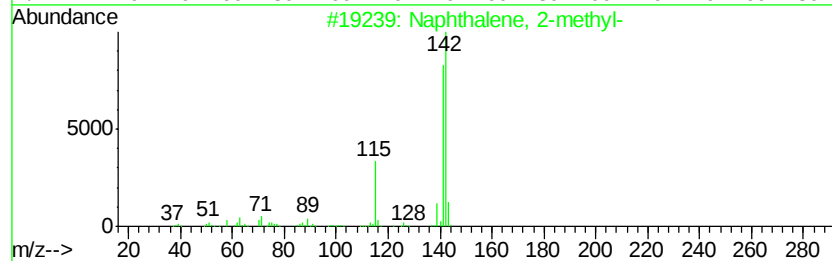
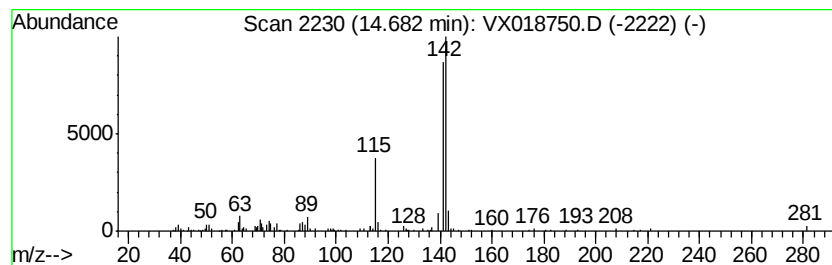
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 Naphthalene, 1-methyl- Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	0.56 ug/L	78567	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100220\
 Data File : VX018750.D
 Acq On : 02 Oct 2020 16:16
 Operator : JC/SP
 Sample : L4171-13DL2 50X
 Misc : 25.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3W0DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
 Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	2.82	2.6	ug/L	226885	1	6.83	436103	5.0
(DEL) Alkane: Str...	2.87	18.0	ug/L	1569380	1	6.83	436103	5.0
(DEL) Alkane: Str...	3.15	43.9	ug/L	3829950	1	6.83	436103	5.0
(DEL) Alkane: Str...	3.50	53.7	ug/L	4686680	1	6.83	436103	5.0
(DEL) Alkane: Str...	4.11	1.6	ug/L	137454	1	6.83	436103	5.0
(DEL) Alkane: Str...	4.28	1.2	ug/L	104790	1	6.83	436103	5.0
(DEL) Alkane: Cyc...	4.37	3.4	ug/L	296958	1	6.83	436103	5.0
Benzene, propyl-	11.34	8.9	ug/L	1251390	3	12.06	703844	5.0
Benzene, 1-ethyl-...	11.42	19.5	ug/L	2739800	3	12.06	703844	5.0
Benzene, 1,2,3-tr...	11.49	11.9	ug/L	1682360	3	12.06	703844	5.0
Benzene, 1-ethyl-...	11.65	2.4	ug/L	342685	3	12.06	703844	5.0
Mesitylene	11.79	17.8	ug/L	2501380	3	12.06	703844	5.0
Benzene, (2-methy...	11.90	1.1	ug/L	157646	3	12.06	703844	5.0
Oxirane, 2-methyl...	11.93	1.3	ug/L	185471	3	12.06	703844	5.0
Benzene, 1-methyl...	12.01	3.0	ug/L	424472	3	12.06	703844	5.0
Benzene, 1,2,4-tr...	12.13	1.4	ug/L	199354	3	12.06	703844	5.0
Benzene, 1,4-diet...	12.29	5.2	ug/L	727662	3	12.06	703844	5.0
Benzene, 1-methyl...	12.30	8.5	ug/L	1194350	3	12.06	703844	5.0
Benzene, 1,2-diet...	12.44	1.2	ug/L	171526	3	12.06	703844	5.0
Benzeneacetaldehy...	12.50	4.3	ug/L	610350	3	12.06	703844	5.0
Benzene, 2-ethyl-...	12.57	13.7	ug/L	1928740	3	12.06	703844	5.0
Benzene, 1,2,4,5-...	12.59	8.5	ug/L	1196200	3	12.06	703844	5.0
Benzene, 2-ethyl-...	12.65	20.8	ug/L	2921590	3	12.06	703844	5.0
o-Cymene	12.75	2.9	ug/L	406767	3	12.06	703844	5.0
Benzene, 4-ethyl-...	12.89	5.5	ug/L	768565	3	12.06	703844	5.0
Benzene, 1,2,3,5-...	12.96	14.4	ug/L	2021330	3	12.06	703844	5.0
Benzene, 1,2,3,4-...	13.00	21.5	ug/L	3023810	3	12.06	703844	5.0
Benzene, 1-methyl...	13.06	0.8	ug/L	107056	3	12.06	703844	5.0
Benzene, 1-etheny...	13.21	4.3	ug/L	604576	3	12.06	703844	5.0
unknown-01	13.29	1.3	ug/L	181541	3	12.06	703844	5.0
unknown-02	13.33	12.9	ug/L	1814540	3	12.06	703844	5.0
Benzene, (1,1-dim...	13.41	1.2	ug/L	174212	3	12.06	703844	5.0
3,4-Dimethylcumene	13.72	1.2	ug/L	168860	3	12.06	703844	5.0
1H-Indene, 2,3-di...	14.12	0.7	ug/L	98188	3	12.06	703844	5.0
1H-Indene, 2,3-di...	14.26	0.6	ug/L	76843	3	12.06	703844	5.0
Naphthalene, 1-me...	14.68	0.6	ug/L	78567	3	12.06	703844	5.0