

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081619\
 Data File : VX011583.D
 Acq On : 16 Aug 2019 16:01
 Operator : JC/SP
 Sample : K4368-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PH2-0133

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 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\82X081519W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.653	731	738	757	rVB	427832	1273909	1.68%	0.508%
2	6.854	925	935	944	rBV	595542	1369746	1.81%	0.546%
3	7.457	1023	1034	1045	rBV	547810	1216201	1.61%	0.485%
4	8.713	1235	1240	1247	rVB	1216678	1865150	2.46%	0.744%
5	10.109	1464	1469	1476	rBV	1166343	1553027	2.05%	0.619%
6	10.250	1486	1492	1501	rBV	16007084	22137261	29.23%	8.827%
7	10.366	1501	1511	1516	rBV	47309289	75735772	100.00%	30.200%
8	11.018	1612	1618	1631	rVB	2645996	3321028	4.39%	1.324%
9	11.134	1631	1637	1642	rBV	1040096	1314237	1.74%	0.524%
10	11.359	1669	1674	1680	rVV	3110745	3850929	5.08%	1.536%
11	11.432	1681	1686	1693	rVV	13637333	22304415	29.45%	8.894%
12	11.506	1693	1698	1707	rVB	7920264	9485463	12.52%	3.782%
13	11.664	1716	1724	1734	rBV	10154883	13073919	17.26%	5.213%
14	11.810	1742	1748	1753	rVB	23789687	30898070	40.80%	12.321%
15	12.024	1778	1783	1786	rBV	710372	855802	1.13%	0.341%
16	12.073	1786	1791	1797	rVB2	1383368	2016092	2.66%	0.804%
17	12.140	1797	1802	1808	rVB	12739865	16278156	21.49%	6.491%
18	12.298	1822	1828	1835	rBV2	4005683	5955343	7.86%	2.375%
19	12.371	1835	1840	1849	rVB2	1711179	2531601	3.34%	1.009%
20	12.518	1859	1864	1869	rBV	1019597	1252246	1.65%	0.499%
21	12.585	1870	1875	1877	rBV	1143253	1539429	2.03%	0.614%
22	12.609	1877	1879	1883	rVV	1270648	1267199	1.67%	0.505%
23	12.664	1883	1888	1891	rVV	1369185	1589142	2.10%	0.634%
24	12.755	1898	1903	1911	rBV2	1145400	1929725	2.55%	0.769%
25	12.902	1922	1927	1931	rVB2	1004043	1261860	1.67%	0.503%
26	12.975	1931	1939	1942	rVV	1128077	1384660	1.83%	0.552%
27	13.017	1942	1946	1952	rVB	1817406	2149366	2.84%	0.857%
28	13.341	1995	1999	2005	rVB2	3273145	4364028	5.76%	1.740%
29	13.475	2017	2021	2027	rVB	1665348	2132657	2.82%	0.850%
30	13.834	2072	2080	2086	rBV	6940270	9276108	12.25%	3.699%
31	14.694	2214	2221	2226	rBV	2541753	3225709	4.26%	1.286%
32	14.834	2239	2244	2249	rBV	1852049	2372144	3.13%	0.946%

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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PH2-0133

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Title : SW846 8260

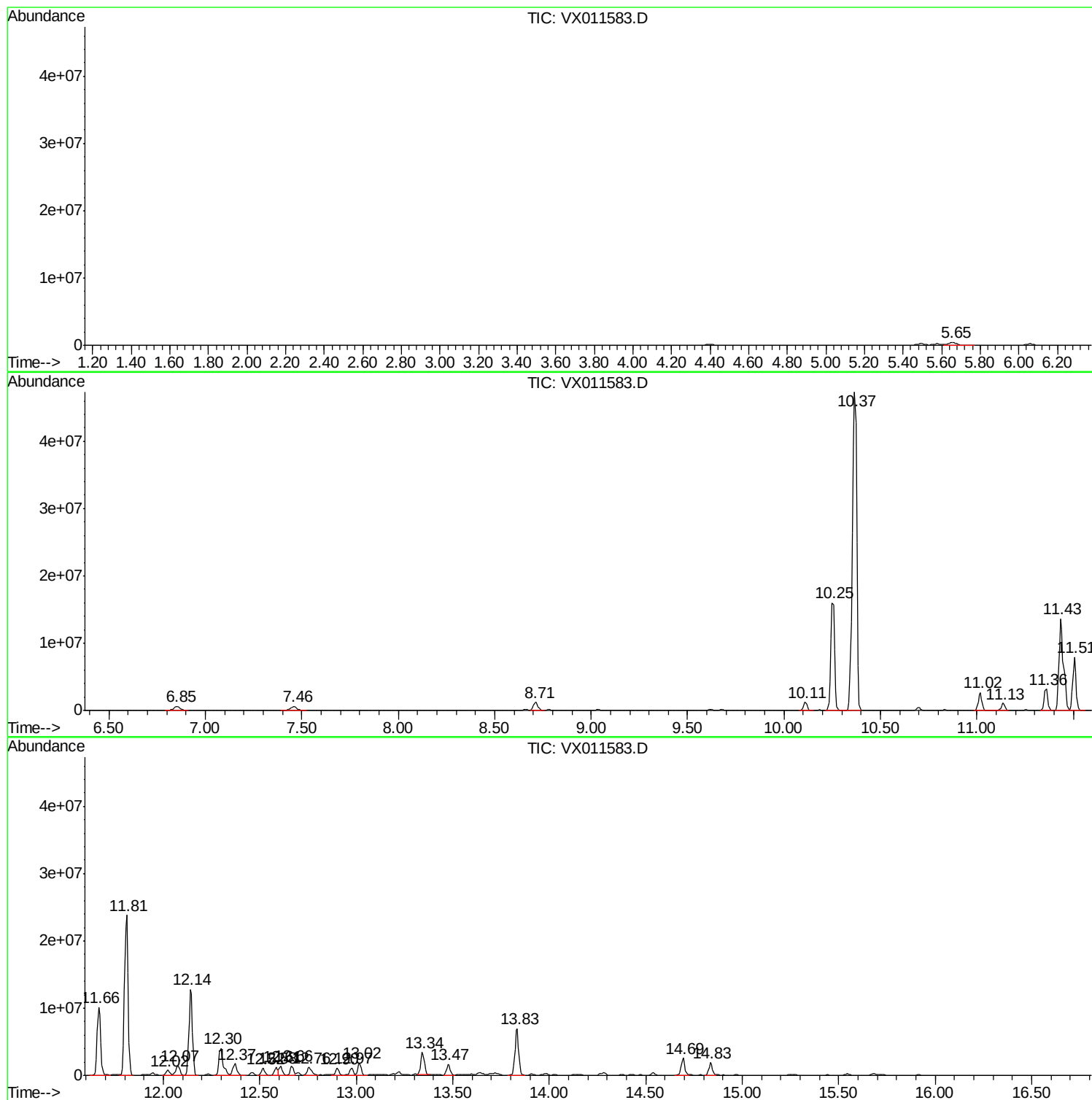
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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



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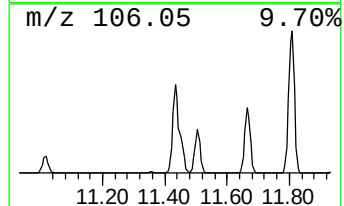
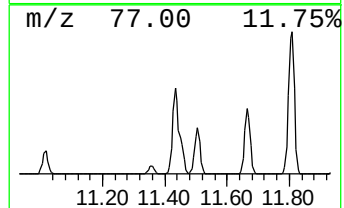
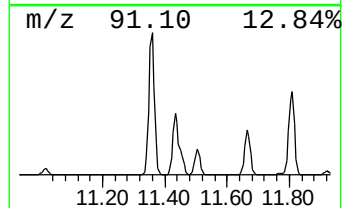
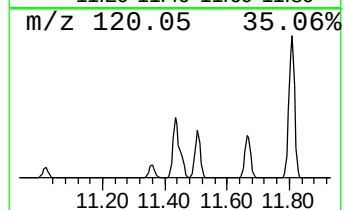
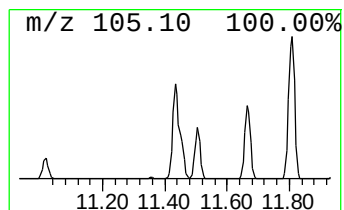
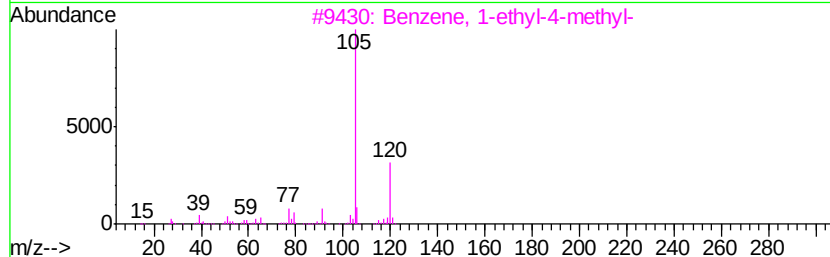
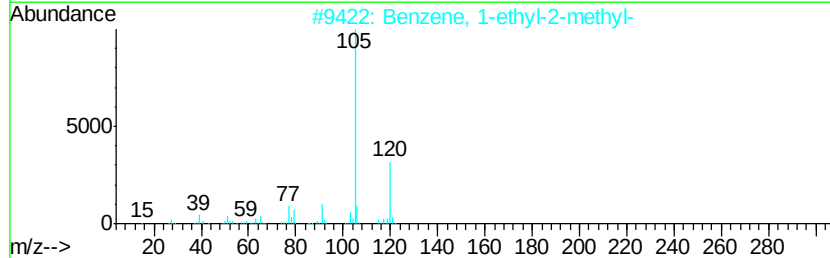
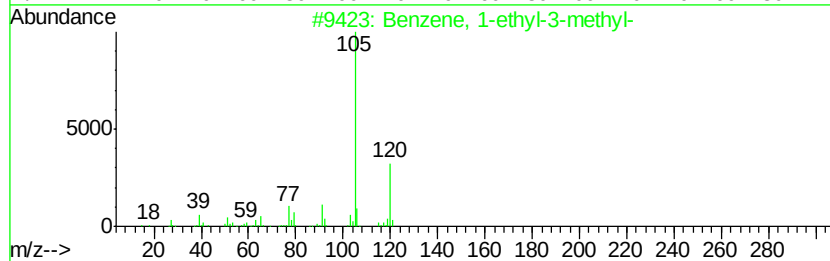
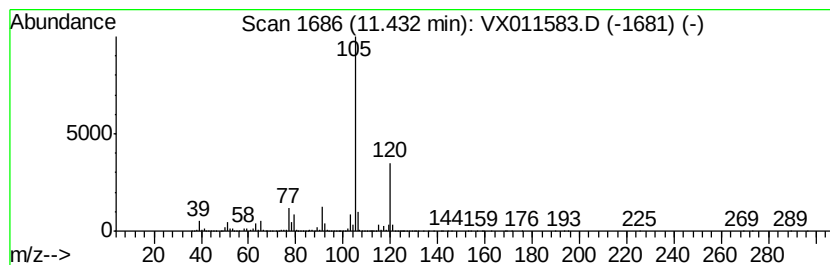
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	553.16 ug/l	22304400	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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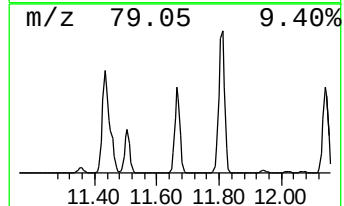
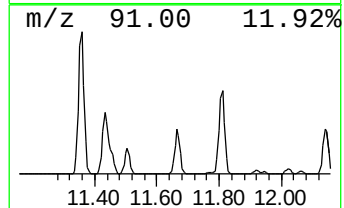
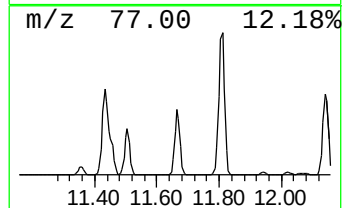
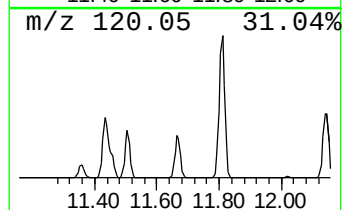
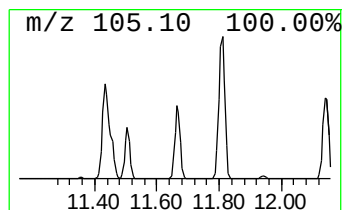
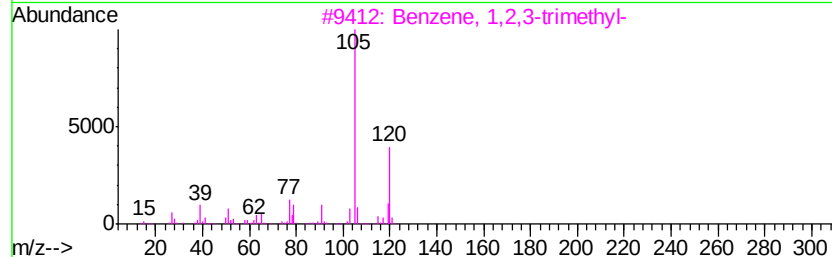
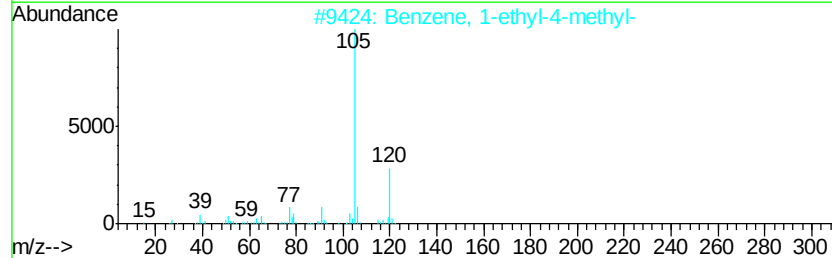
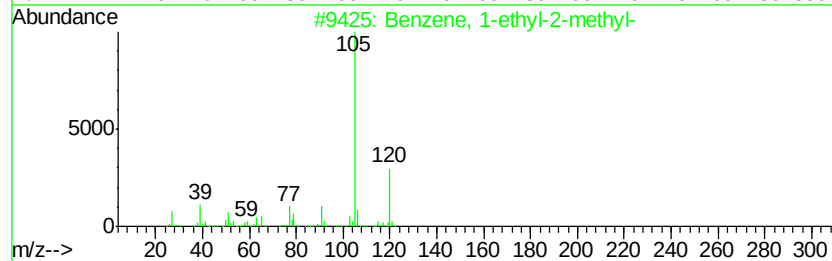
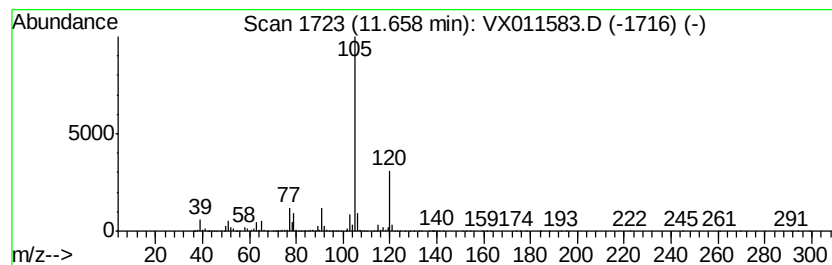
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	324.24 ug/l	13073900	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Mesitylene	120	C9H12	000108-67-8	91



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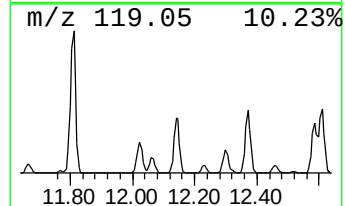
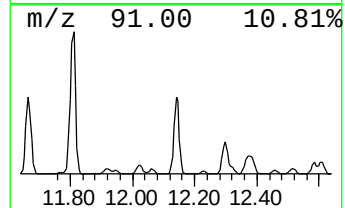
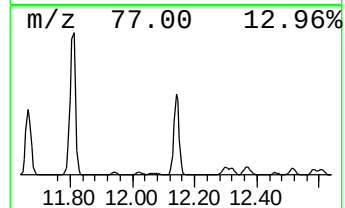
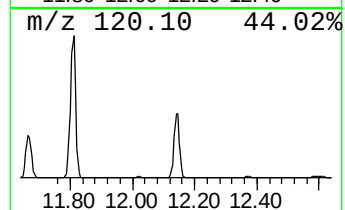
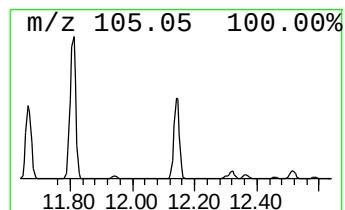
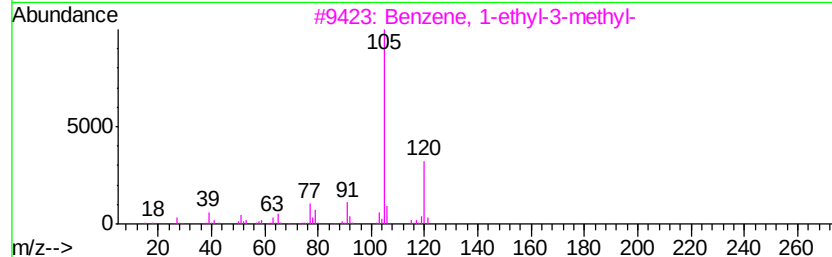
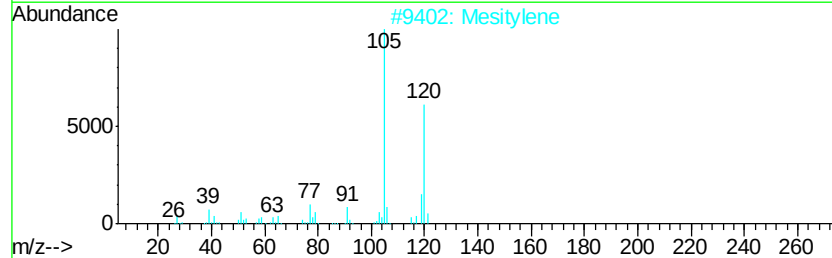
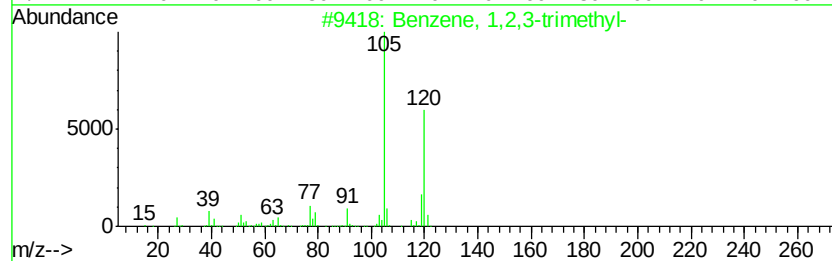
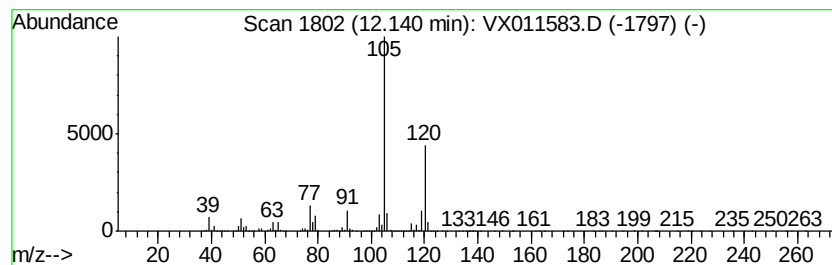
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	403.71 ug/l	16278200	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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Sample : K4368-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

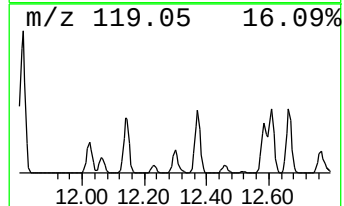
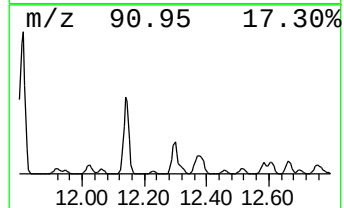
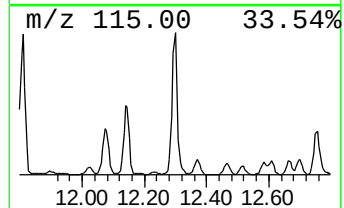
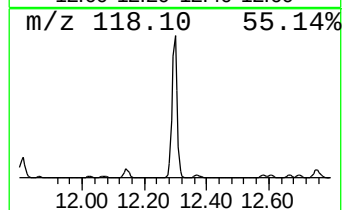
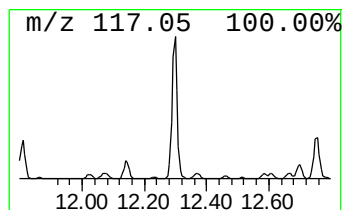
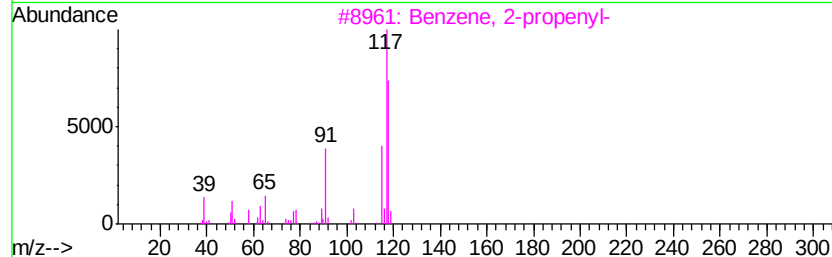
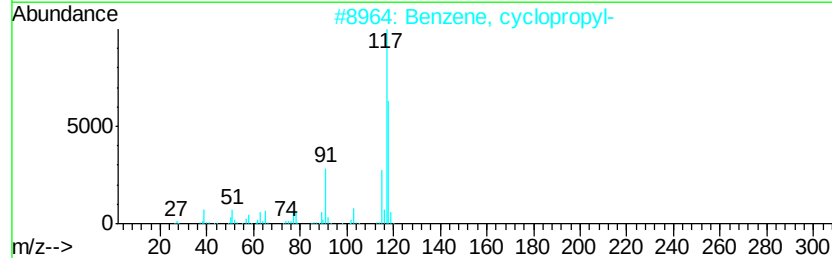
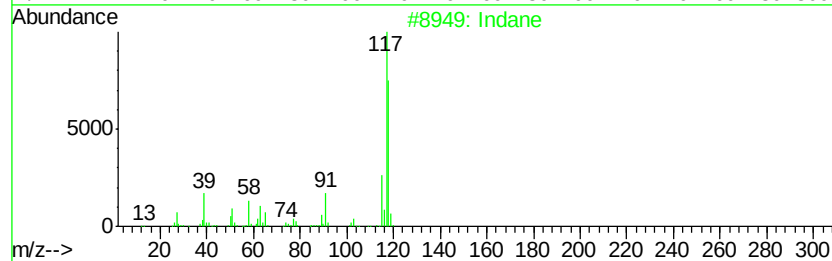
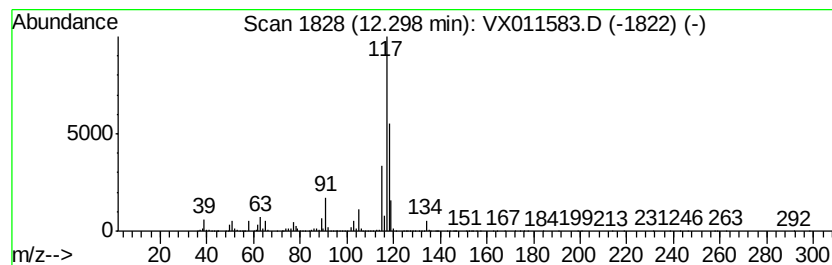
Instrument :
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ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	147.70 ug/l	5955340	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 76
2	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76
3	Benzene, 2-propenyl-		118 C9H10	000300-57-2 68
4	Deltacyclene		118 C9H10	007785-10-6 58
5	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 55



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Instrument :
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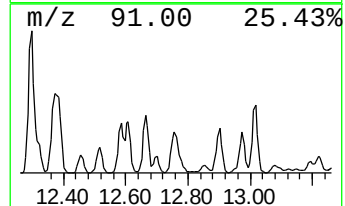
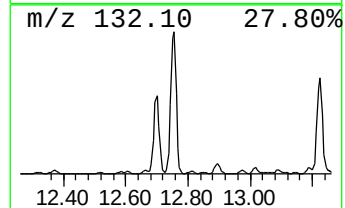
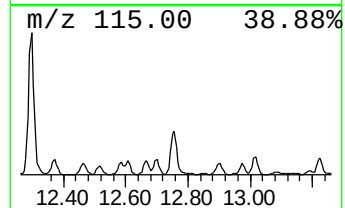
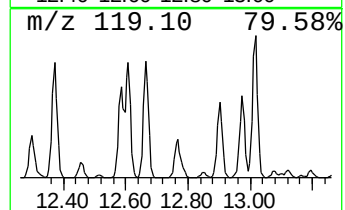
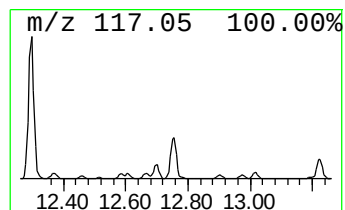
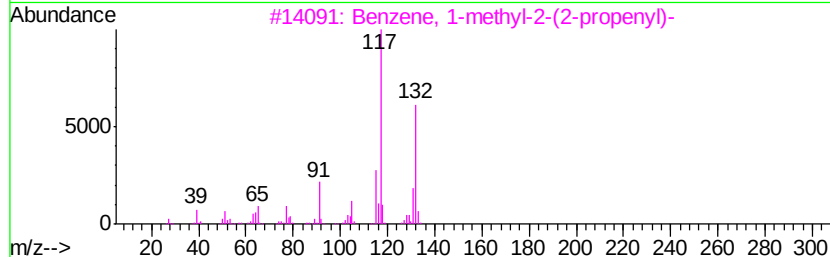
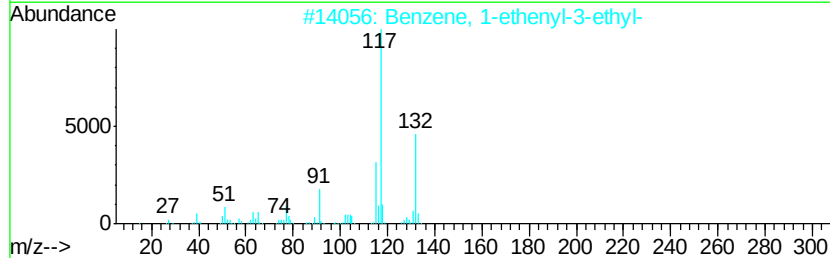
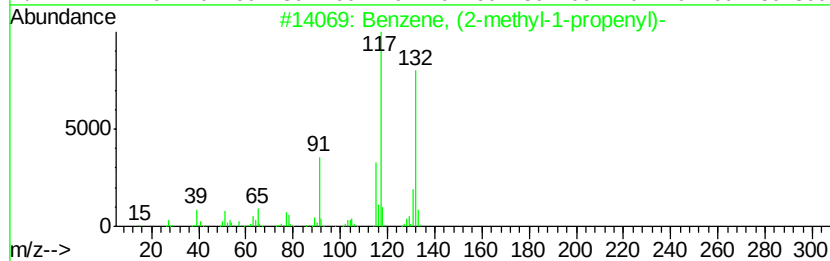
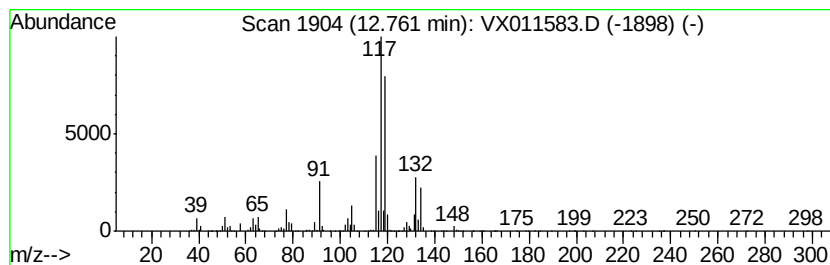
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-propen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	47.86 ug/l	1929730	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
4		Indan, 1-methyl-	132	C10H12	000767-58-8	76
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



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ClientSampleId :
PH2-0133

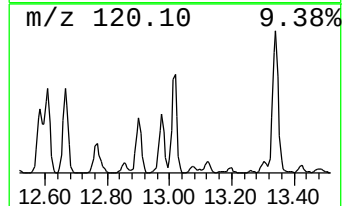
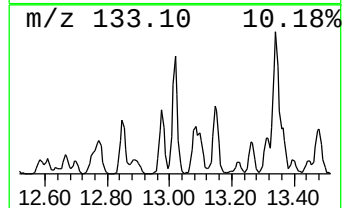
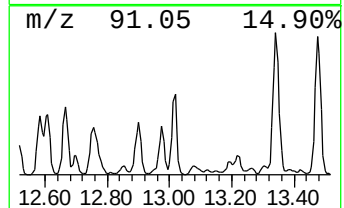
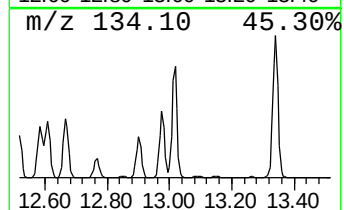
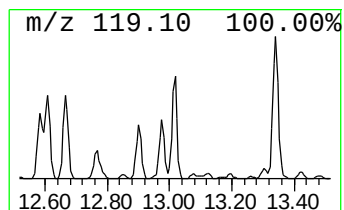
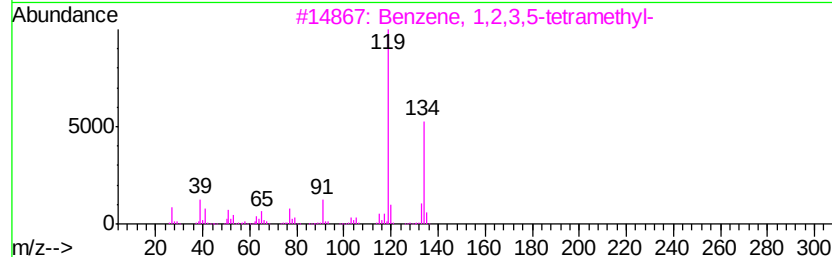
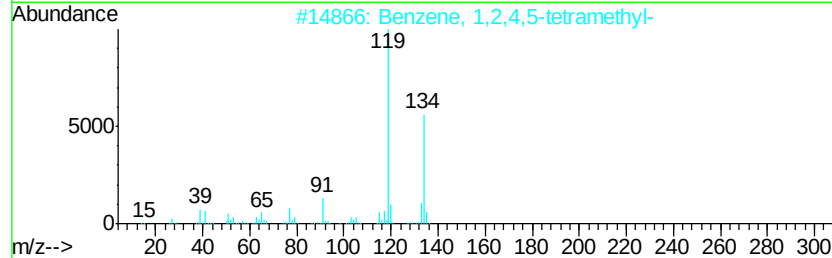
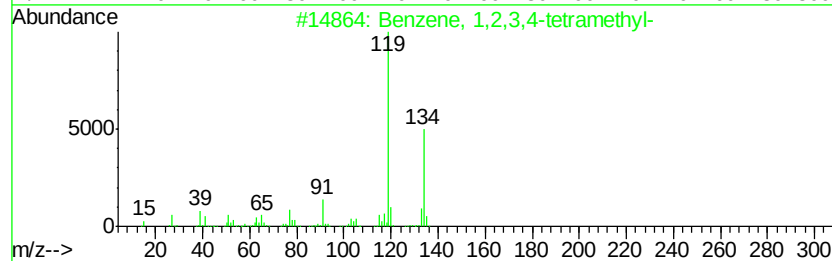
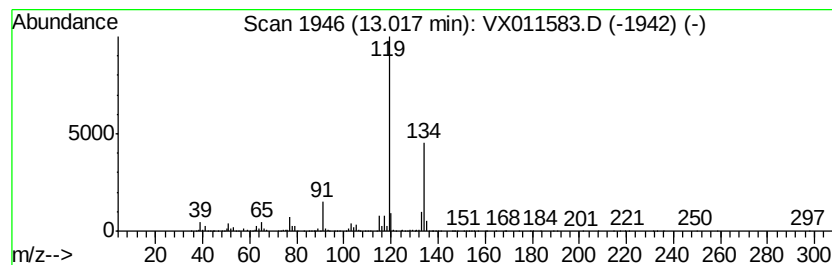
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	53.31 ug/l	2149370	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081619\
Data File : VX011583.D
Acq On : 16 Aug 2019 16:01
Operator : JC/SP
Sample : K4368-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PH2-0133

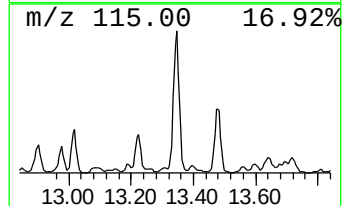
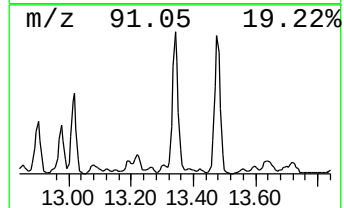
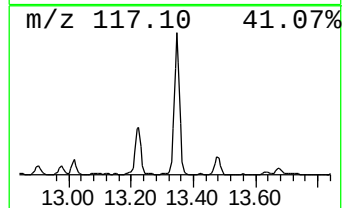
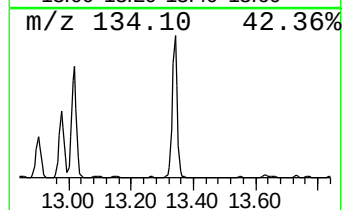
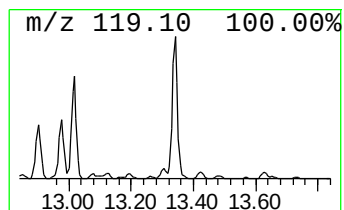
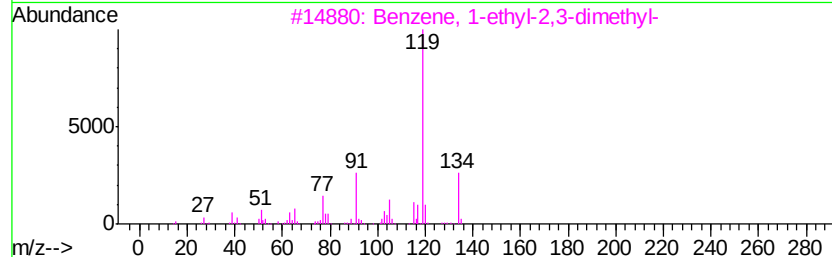
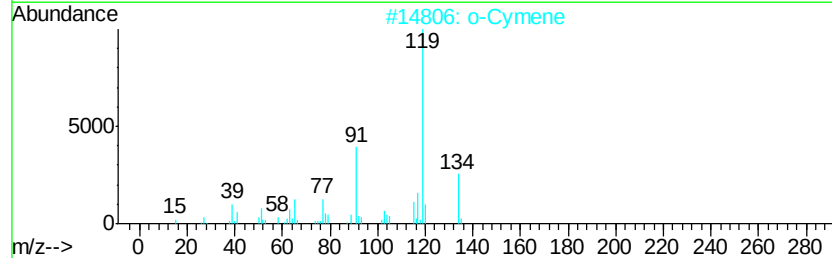
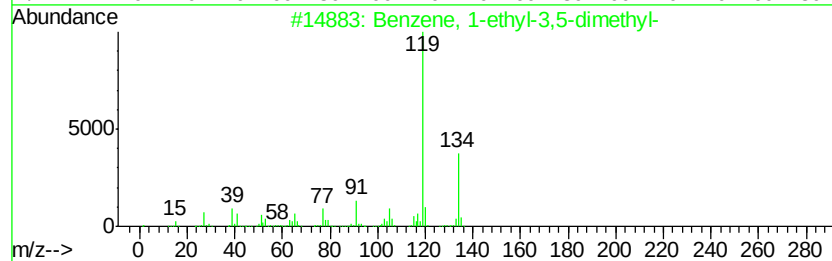
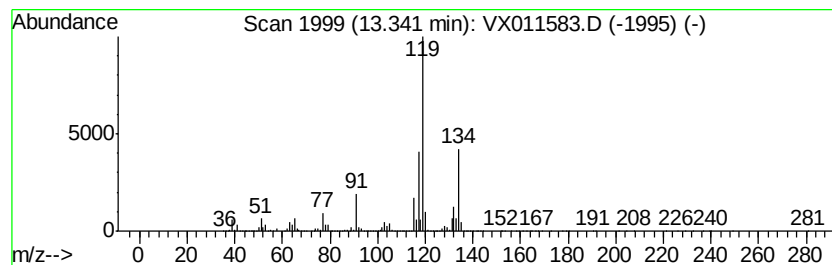
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	108.23 ug/l	4364030	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	76
2		o-Cymene	134	C10H14	000527-84-4	70
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081619\
Data File : VX011583.D
Acq On : 16 Aug 2019 16:01
Operator : JC/SP
Sample : K4368-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
PH2-0133

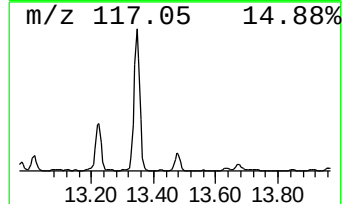
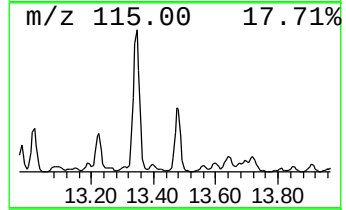
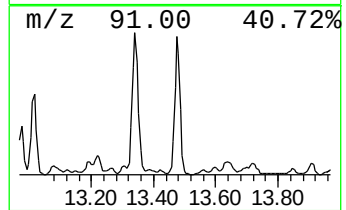
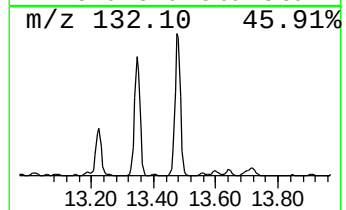
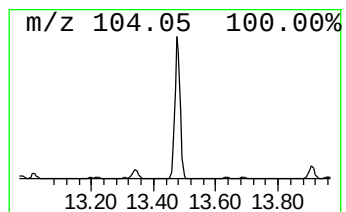
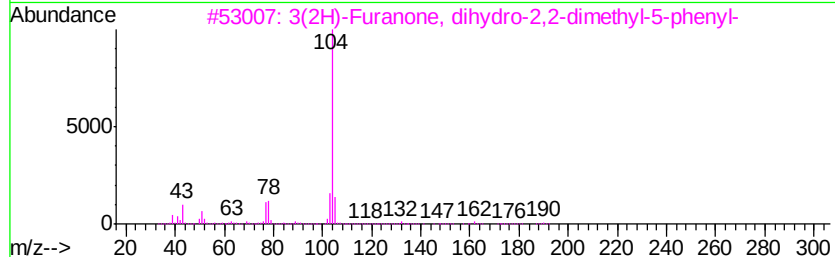
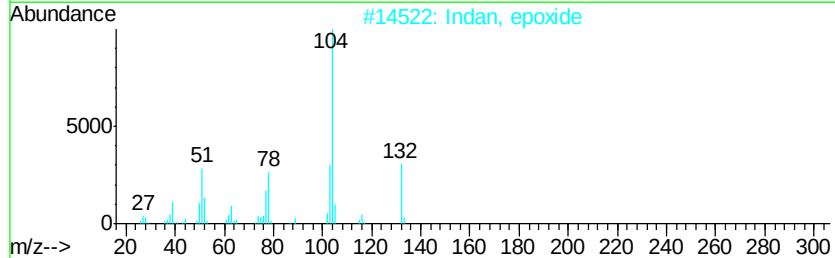
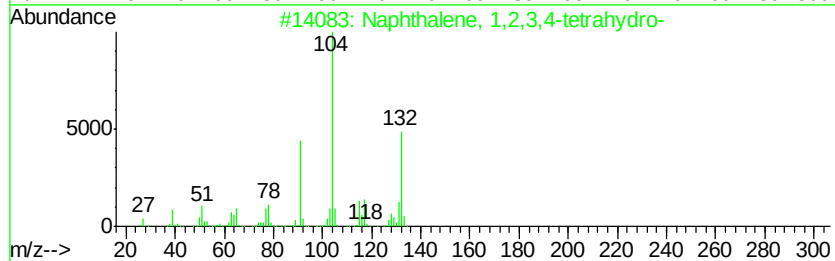
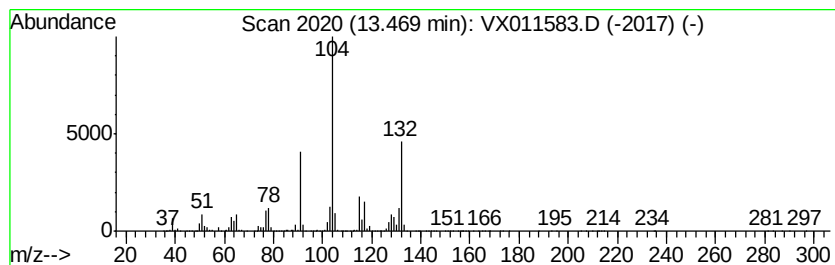
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	52.89 ug/l	2132660	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2		Indan, epoxide	132	C9H8O	000768-22-9	52
3		3(2H)-Furanone, dihydro-2,2-dimethyl-5-phenyl-	190	C12H14O2	063678-00-2	38
4		2-Azetidinone, 4-phenyl-	147	C9H9NO	005661-55-2	38
5		Benzene, 1,1'-(1,2-cyclobutanediyl)-	208	C16H16	007694-30-6	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081619\
Data File : VX011583.D
Acq On : 16 Aug 2019 16:01
Operator : JC/SP
Sample : K4368-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PH2-0133

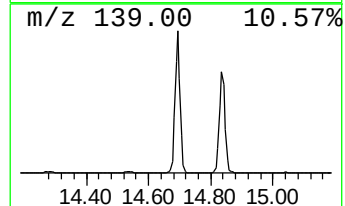
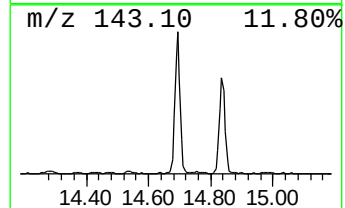
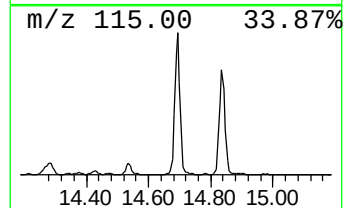
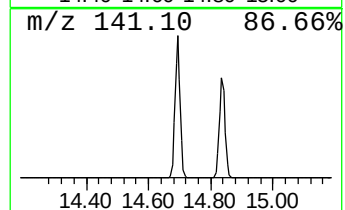
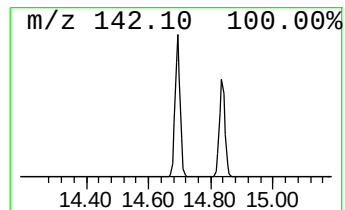
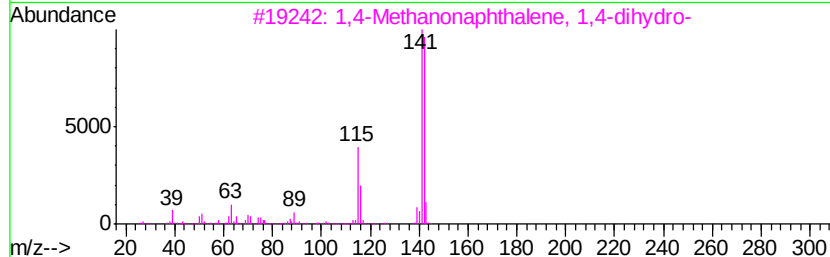
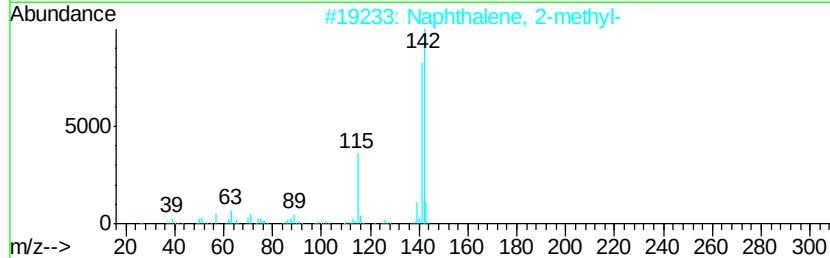
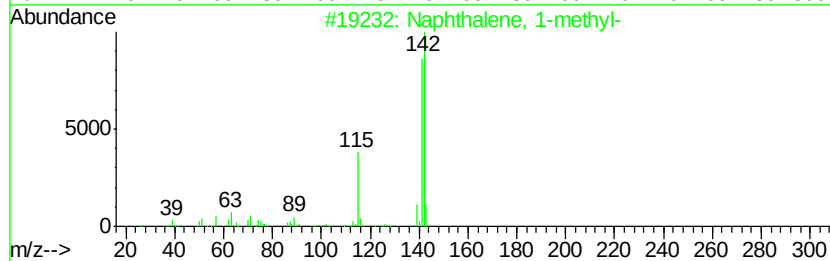
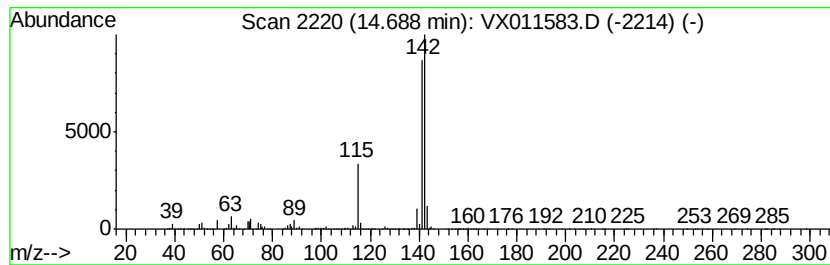
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	80.00 ug/l	3225710	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX081619\
Data File : VX011583.D
Acq On : 16 Aug 2019 16:01
Operator : JC/SP
Sample : K4368-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PH2-0133

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	11.43	553.2	ug/l	22304400	4	12.08	2016090	50.0
Benzene, 1-ethyl-...	11.66	324.2	ug/l	13073900	4	12.08	2016090	50.0
Benzene, 1,2,3-tr...	12.14	403.7	ug/l	16278200	4	12.08	2016090	50.0
Indane	12.30	147.7	ug/l	5955340	4	12.08	2016090	50.0
Benzene, (2-methy...	12.76	47.9	ug/l	1929730	4	12.08	2016090	50.0
Benzene, 1,2,3,4-...	13.02	53.3	ug/l	2149370	4	12.08	2016090	50.0
Benzene, 1-ethyl-...	13.34	108.2	ug/l	4364030	4	12.08	2016090	50.0
Naphthalene, 1,2,...	13.47	52.9	ug/l	2132660	4	12.08	2016090	50.0
Naphthalene, 1-me...	14.69	80.0	ug/l	3225710	4	12.08	2016090	50.0