

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD122818\
 Data File : VD060750.D
 Acq On : 28 Dec 2018 16:59
 Operator : VA/AP
 Sample : J6590-01
 Misc : 5.85g/5ml/MSVOA D/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 AOC2-06A

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_D\METHOD\82D122018S.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	1.439	13	22	47	rBV	1087269	4527921	1.23%	0.127%
2	1.715	47	50	60	rBV	7876496	20593727	5.61%	0.577%
3	1.853	60	64	98	rVB	12737786	64465719	17.57%	1.808%
4	2.354	109	115	133	rBV2	21535444	193644788	52.77%	5.430%
5	2.620	137	142	152	rVB	10501178	51362672	14.00%	1.440%
6	2.945	170	175	189	rVB	10008847	46231149	12.60%	1.296%
7	3.142	189	195	234	rVB2	5148621	30515783	8.32%	0.856%
8	3.801	235	262	287	rBV4	25018179	366994227	100.00%	10.291%
9	4.165	287	299	315	rVB	19714858	158441383	43.17%	4.443%
10	4.401	315	323	331	rBV2	10065799	55604940	15.15%	1.559%
11	4.578	331	341	349	rBV2	14027239	109100695	29.73%	3.059%
12	5.228	400	407	408	rBV	10117189	33693545	9.18%	0.945%
13	5.435	414	428	430	rBV3	13411398	82317743	22.43%	2.308%
14	5.976	466	483	487	rBV3	3238663	18452272	5.03%	0.517%
15	6.143	488	500	513	rVV3	9960250	78963909	21.52%	2.214%
16	6.547	518	541	543	rBV6	18982265	189704788	51.69%	5.320%
17	7.078	578	595	597	rBV	20908372	146894216	40.03%	4.119%
18	7.403	619	628	634	rVB4	15784973	91128910	24.83%	2.555%
19	7.550	634	643	656	rBV4	15397941	116561105	31.76%	3.269%
20	7.728	656	661	665	rBV	5570803	21151743	5.76%	0.593%
21	7.836	665	672	689	rVB6	16500824	122074824	33.26%	3.423%
22	8.161	689	705	708	rBV5	21367126	163119314	44.45%	4.574%
23	8.534	736	743	744	rBV3	6232098	19182934	5.23%	0.538%
24	8.771	755	767	768	rBV	16383557	82162649	22.39%	2.304%
25	9.548	833	846	851	rBV3	16597265	94175449	25.66%	2.641%
26	9.755	861	867	870	rBV3	11216670	45006460	12.26%	1.262%
27	9.991	887	891	893	rBV2	4594457	12199991	3.32%	0.342%
28	10.198	904	912	915	rBV4	12879525	55897095	15.23%	1.567%
29	10.266	916	919	923	rVV2	12313081	36072695	9.83%	1.012%
30	10.375	927	930	934	rBV	7059379	15018493	4.09%	0.421%
31	10.512	939	944	948	rBV3	11039854	36624932	9.98%	1.027%
32	10.660	951	959	964	rBV2	10228915	47857711	13.04%	1.342%
33	10.837	973	977	980	rBV4	6566588	17413763	4.74%	0.488%
34	10.955	986	989	994	rVB4	6549757	20265612	5.52%	0.568%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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35	11.034	994	997	1000	rBV	6453739	16454630	4.48%	0.461%
36	11.152	1000	1009	1015	rVB3	12858545	65483463	17.84%	1.836%
37	11.280	1015	1022	1028	rVB	14216077	58865791	16.04%	1.651%
38	11.379	1028	1032	1036	rBV3	14168839	46961920	12.80%	1.317%
39	11.487	1040	1043	1046	rBV2	6055774	11812334	3.22%	0.331%
40	11.684	1060	1063	1069	rVB2	18239869	52622505	14.34%	1.476%
41	11.772	1069	1072	1077	rVB2	6299247	18967887	5.17%	0.532%
42	11.861	1077	1081	1083	rBV3	6562847	16541680	4.51%	0.464%
43	11.920	1083	1087	1092	rBV5	9170282	32316753	8.81%	0.906%
44	11.998	1093	1095	1100	rVB	5635357	13010088	3.55%	0.365%
45	12.067	1100	1102	1104	rBV2	3658031	7650599	2.08%	0.215%
46	12.313	1122	1127	1130	rBV	5799091	14862814	4.05%	0.417%
47	12.392	1130	1135	1138	rBV3	5660986	19212141	5.23%	0.539%
48	12.491	1141	1145	1150	rVV	13327828	41409017	11.28%	1.161%
49	12.579	1151	1154	1156	rVV	5973369	11711309	3.19%	0.328%
50	12.668	1160	1163	1167	rVV	9022755	23336970	6.36%	0.654%
51	12.756	1168	1172	1174	rVV	8526970	22915440	6.24%	0.643%
52	12.845	1178	1181	1187	rVB3	16142507	48778257	13.29%	1.368%
53	12.983	1187	1195	1197	rBV2	9692195	34460886	9.39%	0.966%
54	13.140	1206	1211	1215	rVB2	15060744	42289199	11.52%	1.186%
55	13.199	1215	1217	1219	rBV	5606012	10190149	2.78%	0.286%
56	13.297	1224	1227	1232	rVB2	7009199	21993961	5.99%	0.617%
57	13.366	1232	1234	1236	rBV	2581491	5283907	1.44%	0.148%
58	13.475	1243	1245	1247	rBV	4105481	5996540	1.63%	0.168%
59	13.593	1253	1257	1258	rBV	9091869	22897570	6.24%	0.642%
60	13.662	1262	1264	1269	rVV2	17158297	43535784	11.86%	1.221%
61	13.730	1269	1271	1272	rVV	5239142	7628122	2.08%	0.214%
62	13.790	1272	1277	1280	rVV2	8656565	25448264	6.93%	0.714%
63	13.849	1280	1283	1285	rVV	10347849	22440872	6.11%	0.629%
64	13.937	1289	1292	1294	rVV	14587055	29297687	7.98%	0.822%
65	14.016	1297	1300	1303	rVB	14626154	27290908	7.44%	0.765%
66	14.075	1303	1306	1309	rVB3	2794557	5489068	1.50%	0.154%
67	14.124	1309	1311	1314	rBV	9247268	13301635	3.62%	0.373%
68	14.203	1317	1319	1321	rVV	9081271	14925192	4.07%	0.419%
69	14.242	1321	1323	1325	rVB	9056612	12593422	3.43%	0.353%
70	14.282	1325	1327	1334	rVB3	6289641	15631487	4.26%	0.438%
71	14.390	1334	1338	1340	rBV3	2660223	5805442	1.58%	0.163%

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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

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72	14.429	1340	1342	1346	rVB	8998869	12485356	3.40%	0.350%
73	14.528	1349	1352	1356	rVB2	10742372	16826891	4.59%	0.472%

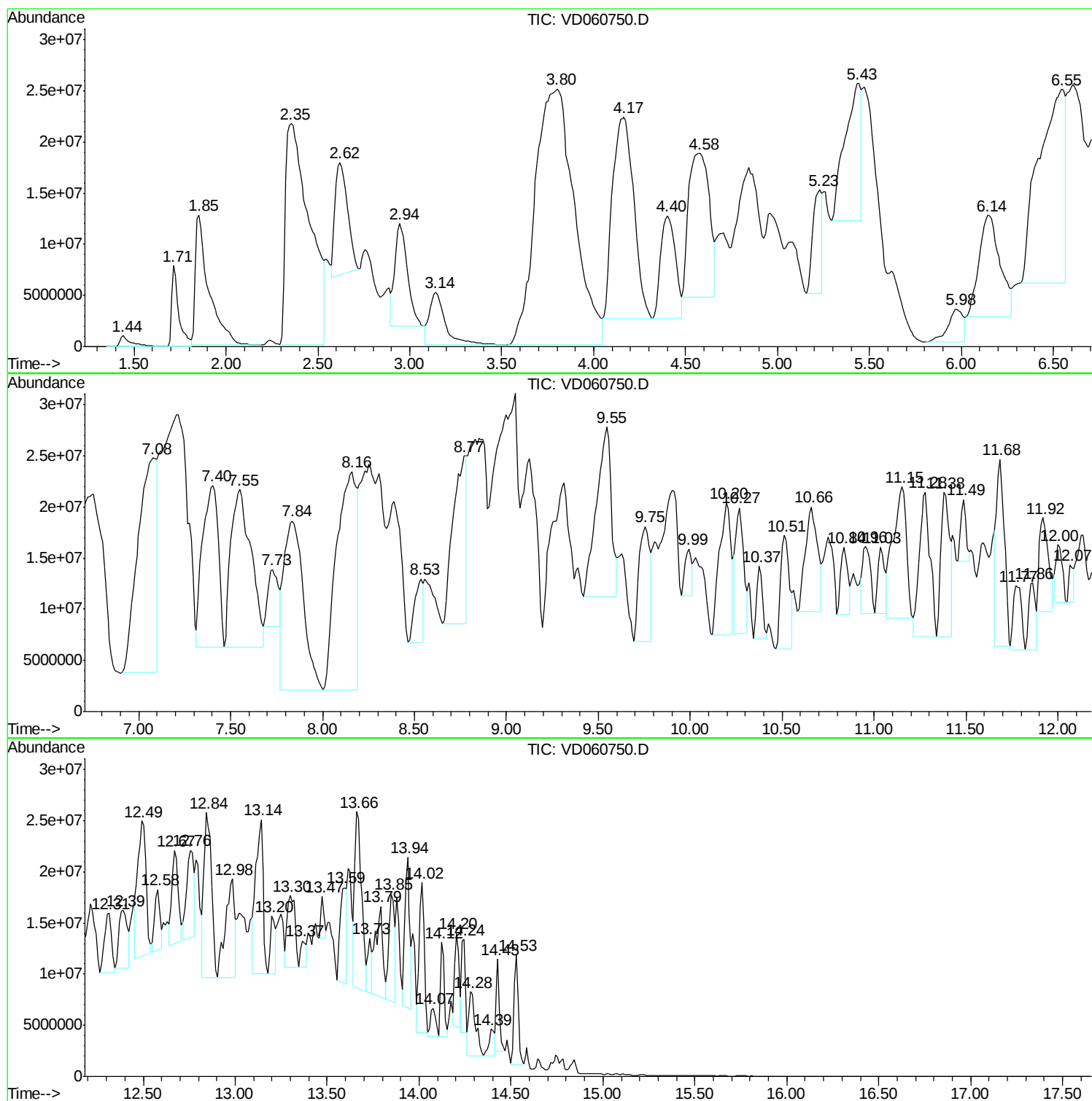
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD122818\
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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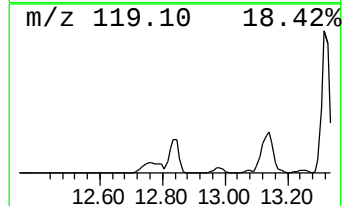
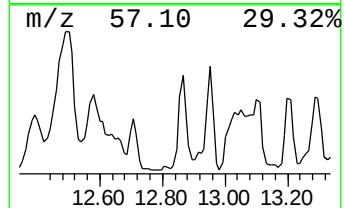
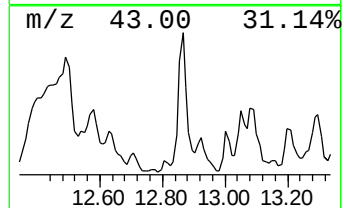
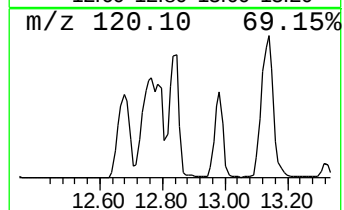
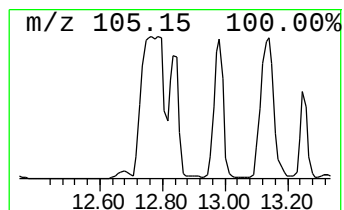
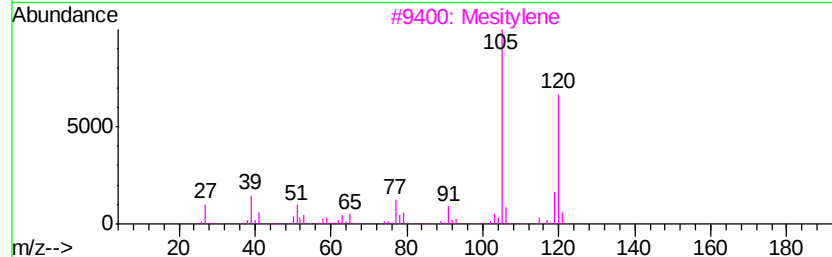
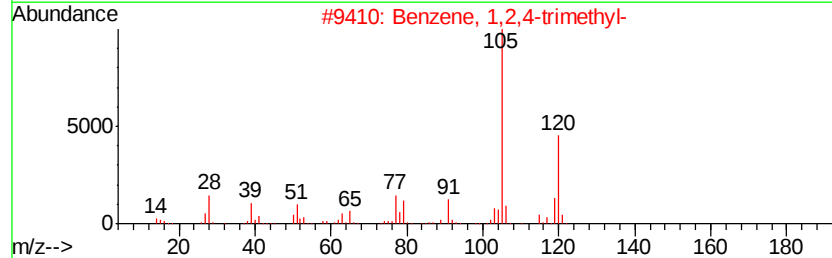
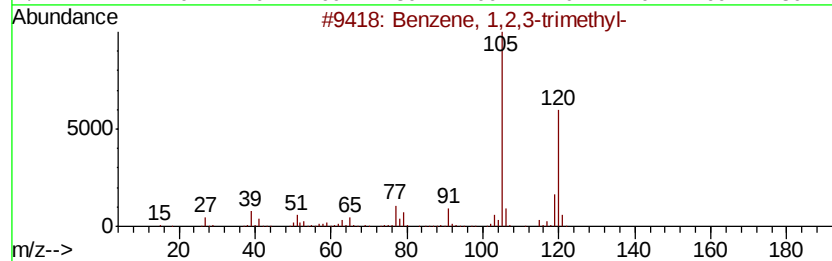
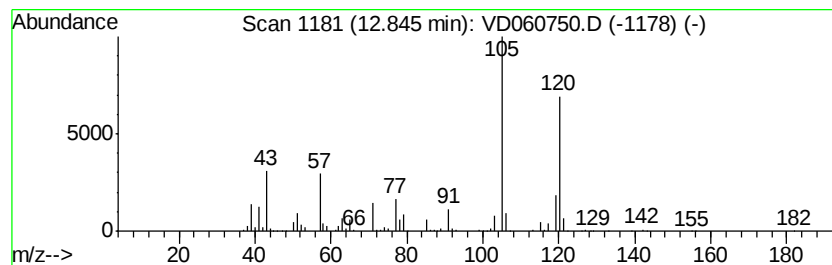
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.84	461.57 ug/l	48778300	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



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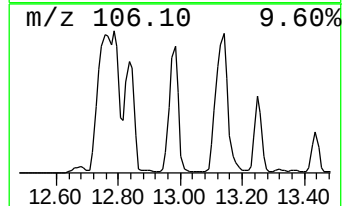
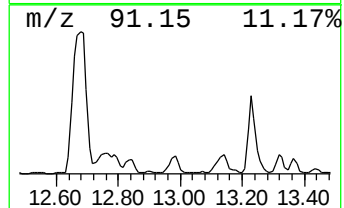
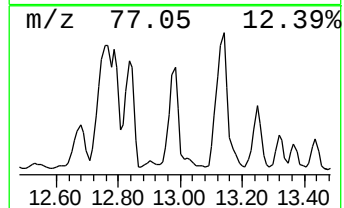
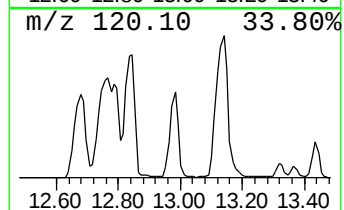
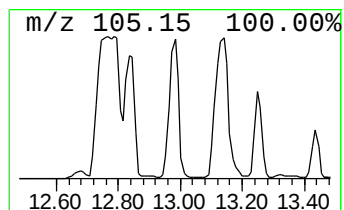
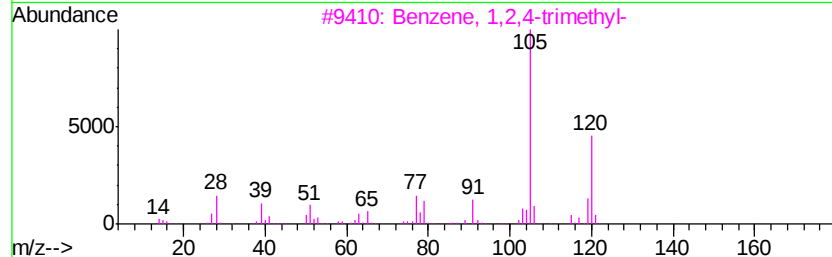
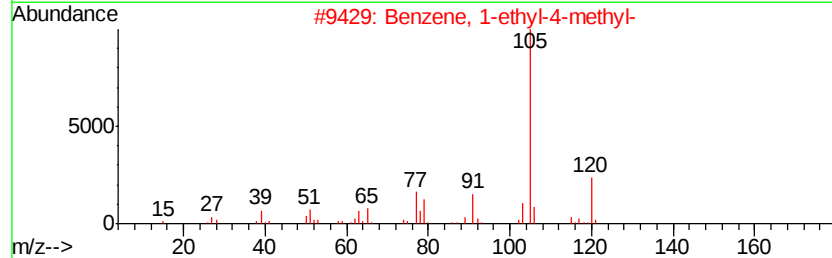
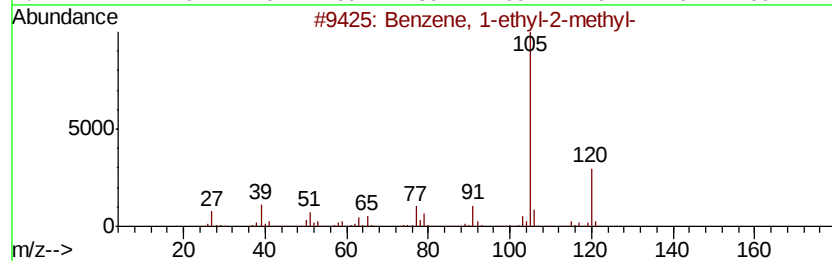
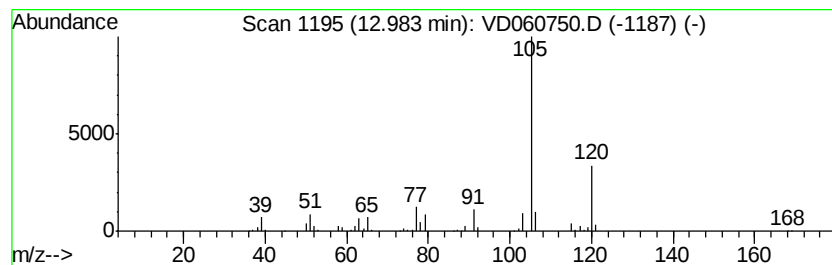
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	326.09 ug/l	34460900	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



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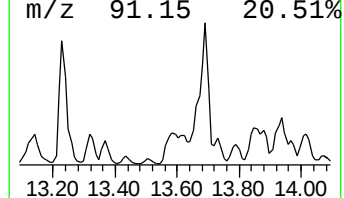
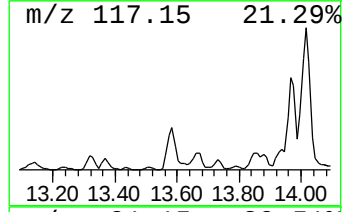
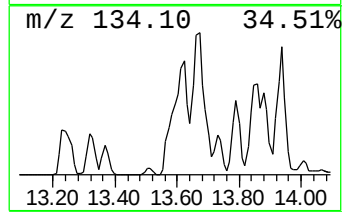
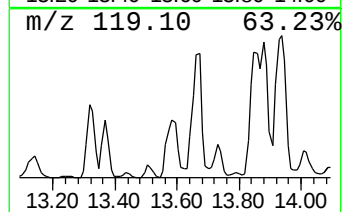
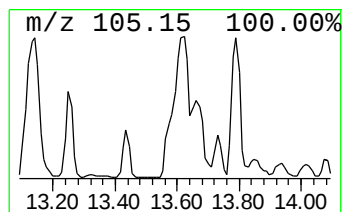
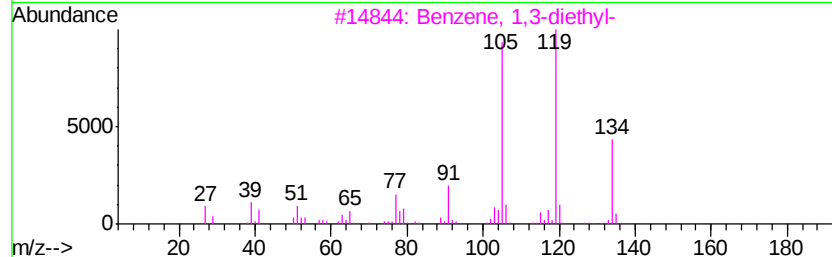
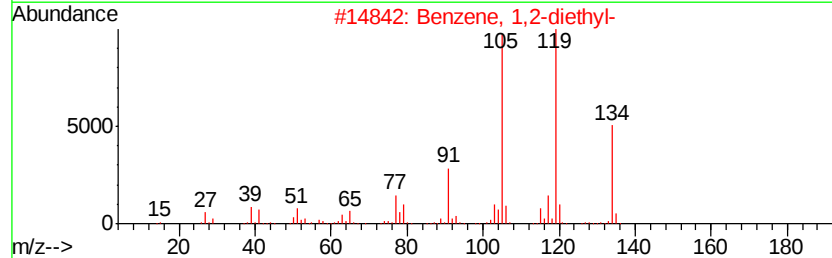
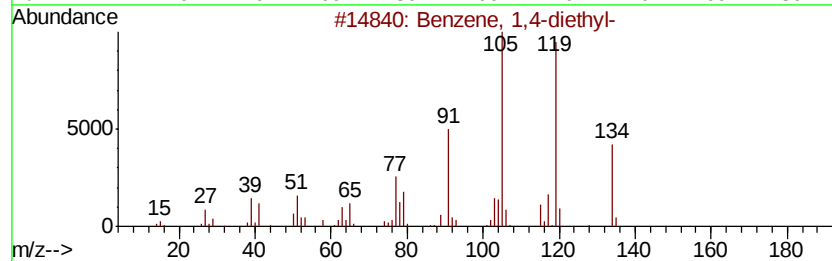
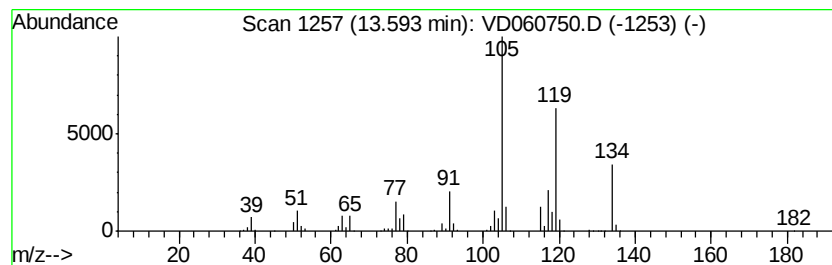
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,4-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.59	216.67 ug/l	22897600	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	64
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	62
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	58
4		2-Indanol	134	C9H10O	004254-29-9	50
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	49



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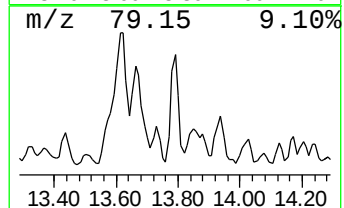
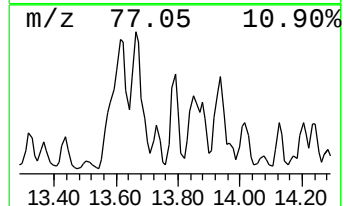
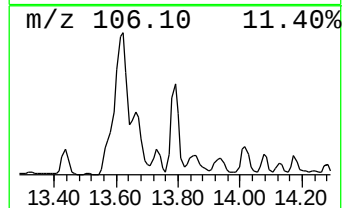
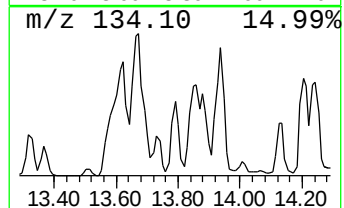
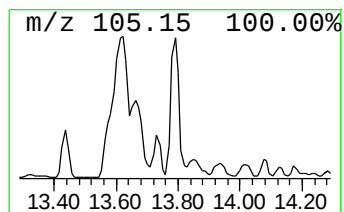
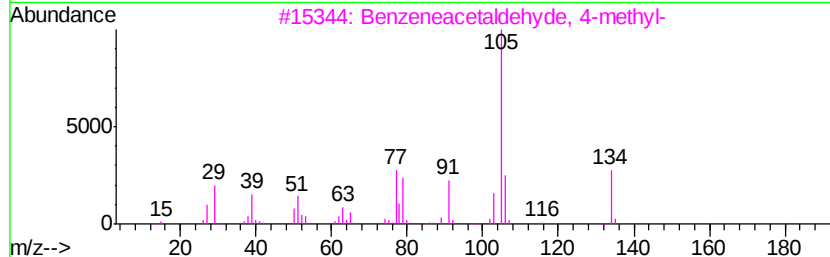
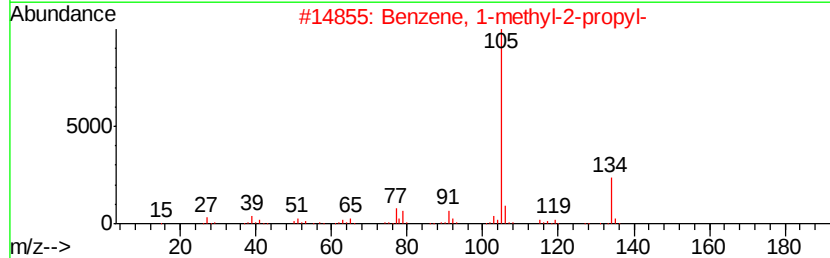
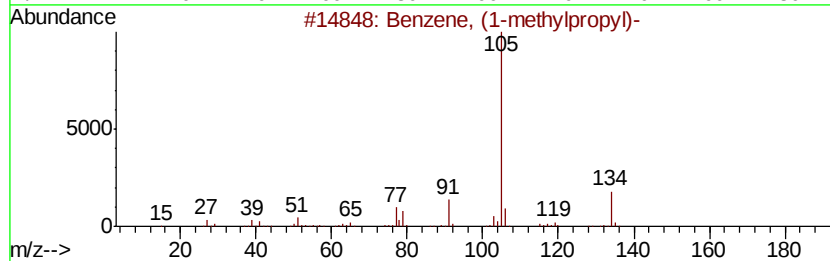
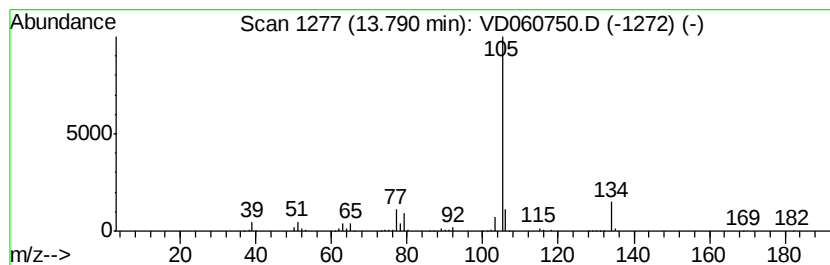
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, (1-methylpropyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	240.81 ug/l	25448300	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	86
3		Benzeneacetaldehyde, 4-methyl-	134	C9H10O	000104-09-6	86
4		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	80
5		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD122818\
Data File : VD060750.D
Acq On : 28 Dec 2018 16:59
Operator : VA/AP
Sample : J6590-01
Misc : 5.85g/5ml/MSVOA D/SOIL
ALS Vial : 9 Sample Multiplier: 1

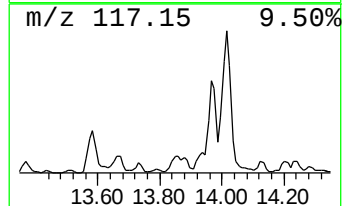
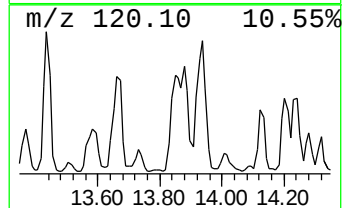
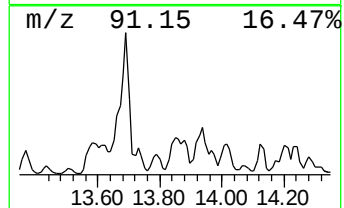
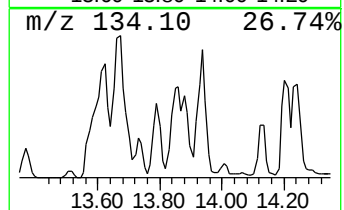
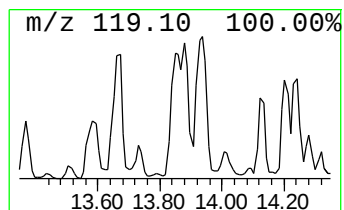
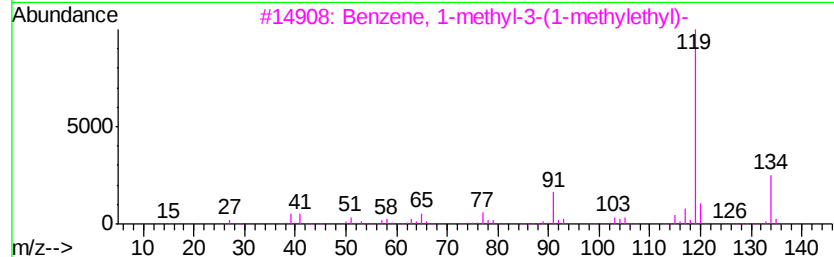
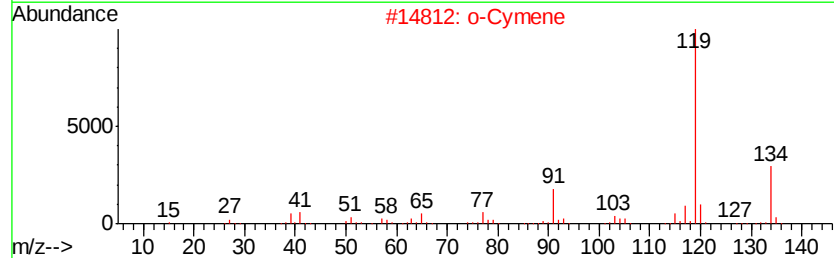
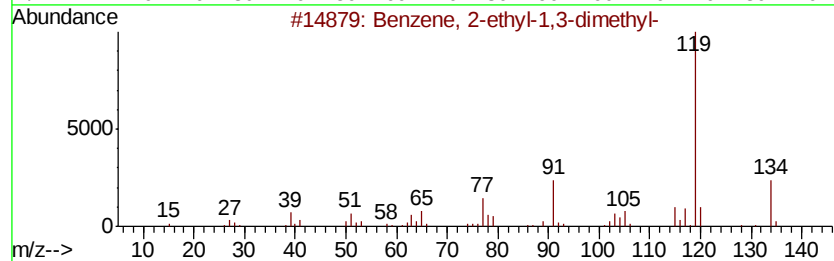
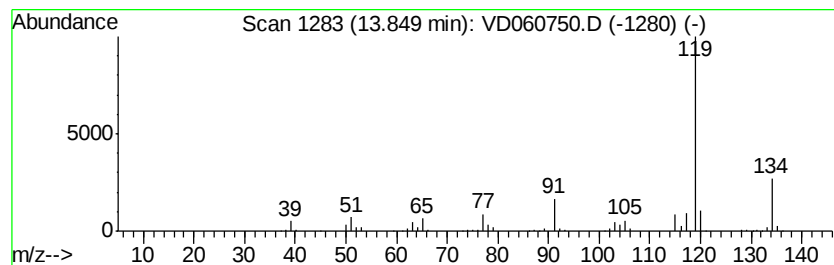
Instrument :
MSVOA_D
ClientSampleId :
AOC2-06A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	212.35 ug/l	22440900	1,4-Dichlorobenzene-d4	13.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	95
2	o-Cymene	134 C10H14	000527-84-4	94
3	Benzene, 1-methyl-3-(1-methylethyl-)	134 C10H14	000535-77-3	94
4	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	91
5	Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD122818\
Data File : VD060750.D
Acq On : 28 Dec 2018 16:59
Operator : VA/AP
Sample : J6590-01
Misc : 5.85g/5ml/MSVOA D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AOC2-06A

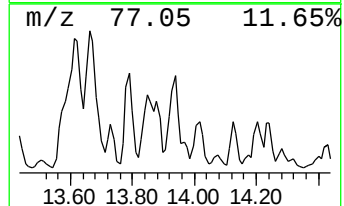
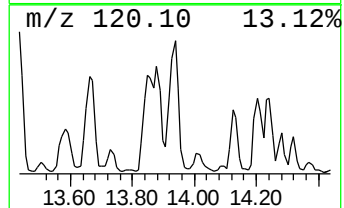
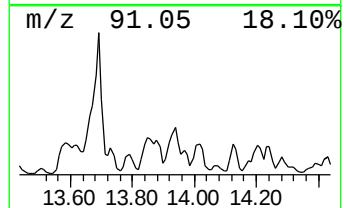
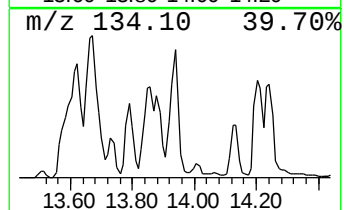
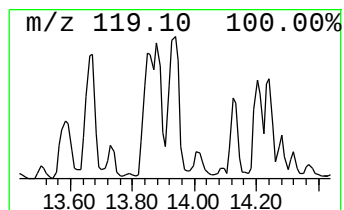
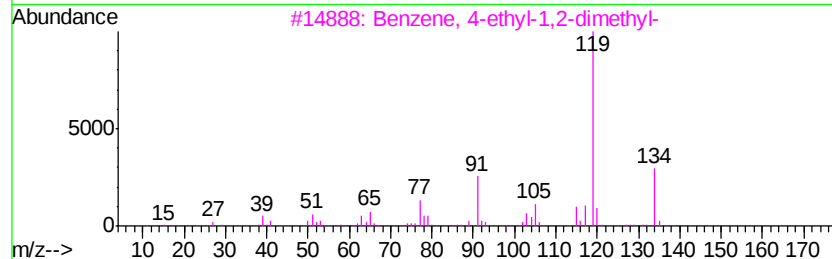
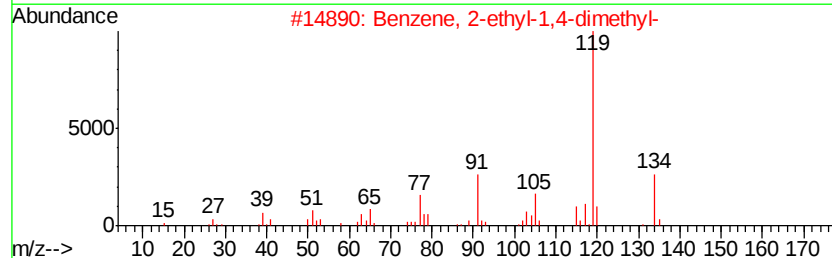
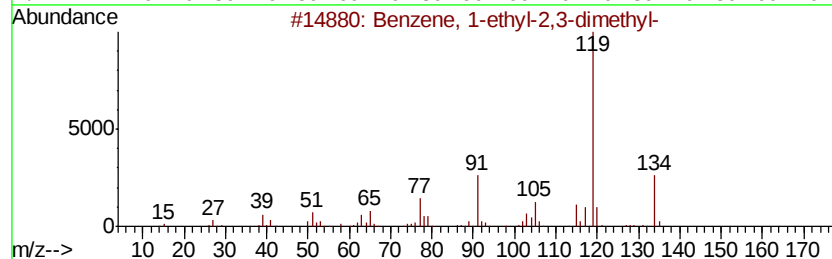
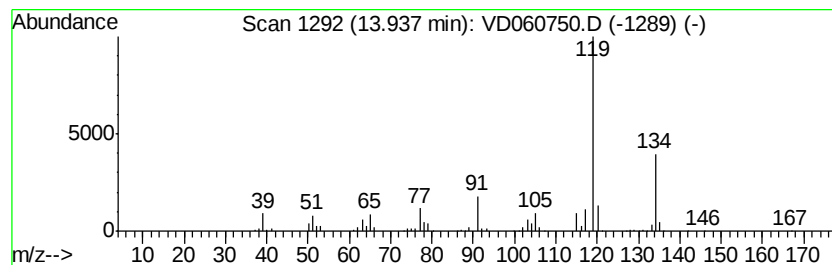
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	277.24 ug/l	29297700	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5		o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD122818\
Data File : VD060750.D
Acq On : 28 Dec 2018 16:59
Operator : VA/AP
Sample : J6590-01
Misc : 5.85g/5ml/MSVOA D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AOC2-06A

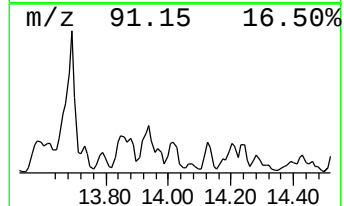
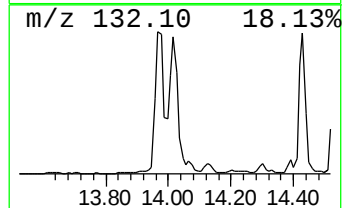
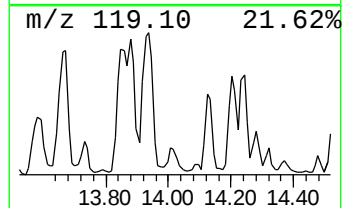
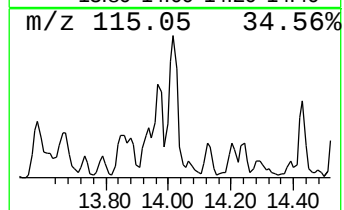
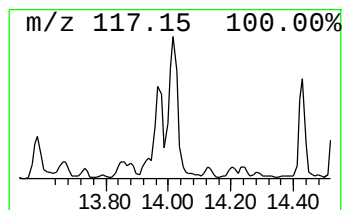
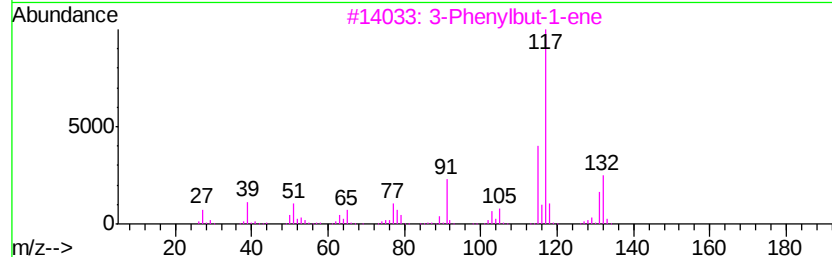
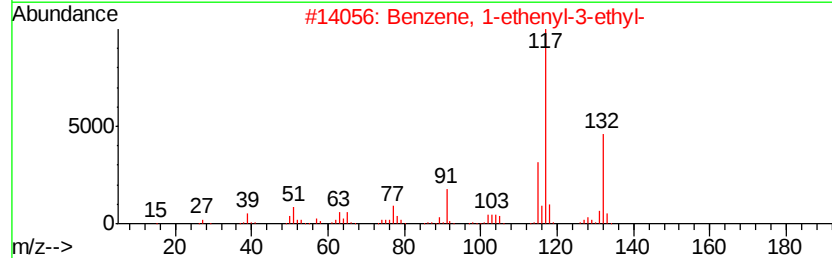
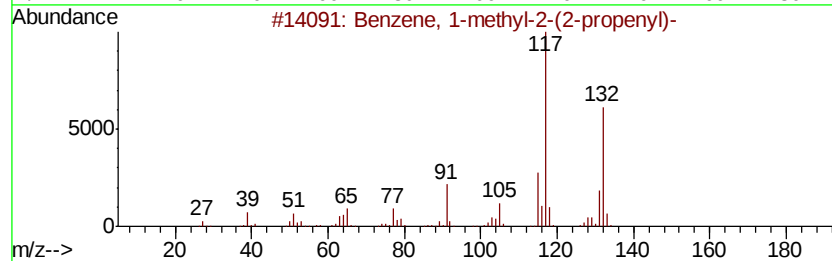
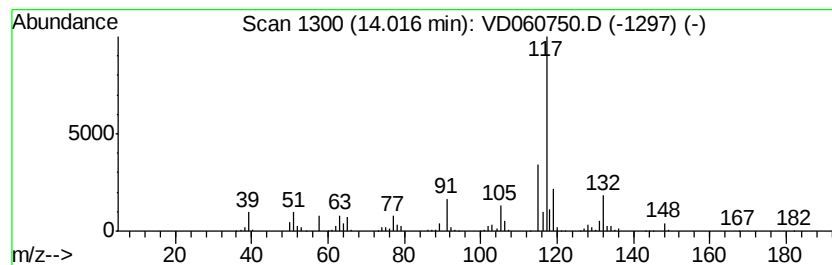
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-2-(2-prop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	258.25 ug/l	27290900	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD122818\
Data File : VD060750.D
Acq On : 28 Dec 2018 16:59
Operator : VA/AP
Sample : J6590-01
Misc : 5.85g/5ml/MSVOA D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AOC2-06A

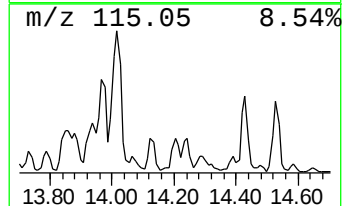
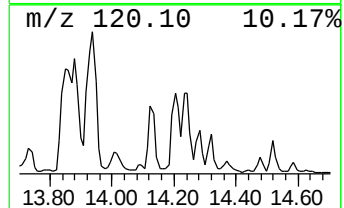
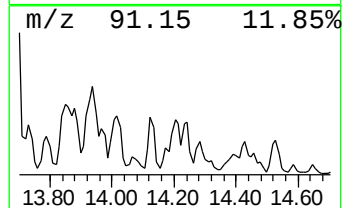
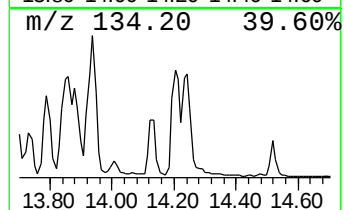
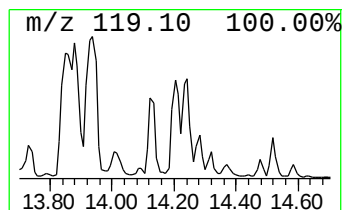
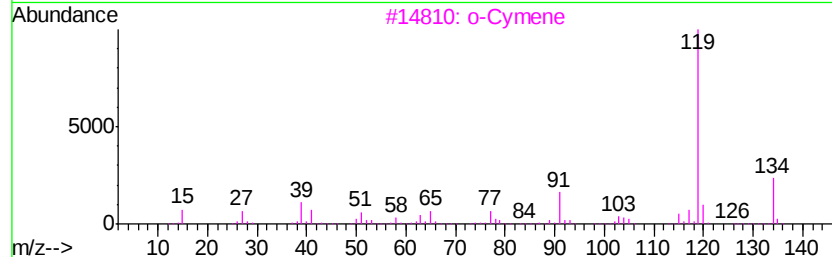
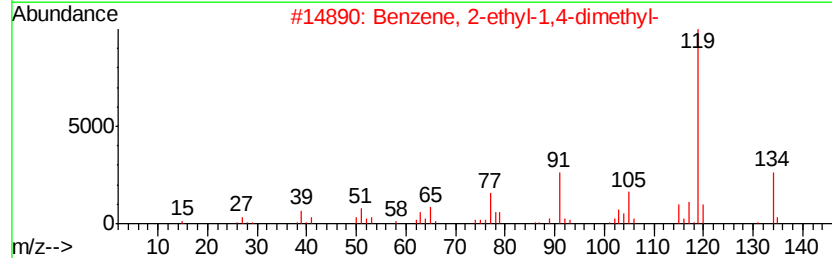
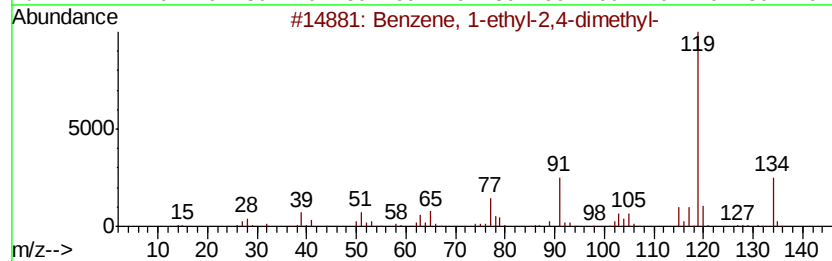
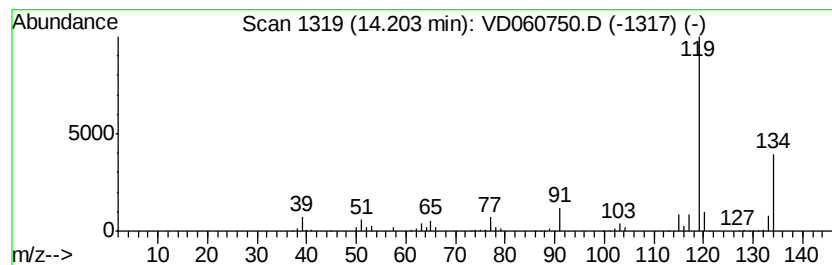
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	141.23 ug/l	14925200	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
3		o-Cymene	134	C10H14	000527-84-4	91
4		p-Cymene	134	C10H14	000099-87-6	91
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD122818\
Data File : VD060750.D
Acq On : 28 Dec 2018 16:59
Operator : VA/AP
Sample : J6590-01
Misc : 5.85g/5ml/MSVOA D/SOIL
ALS Vial : 9 Sample Multiplier: 1

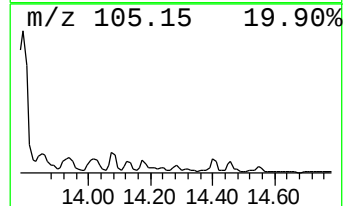
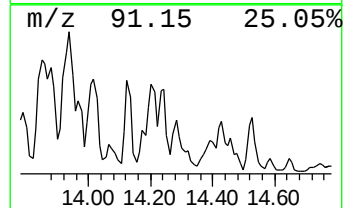
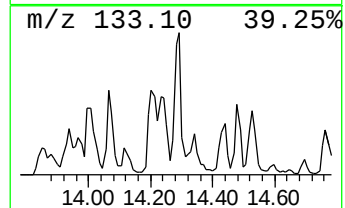
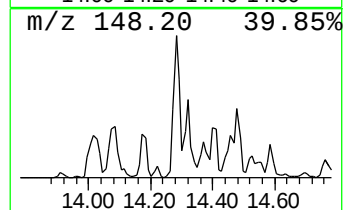
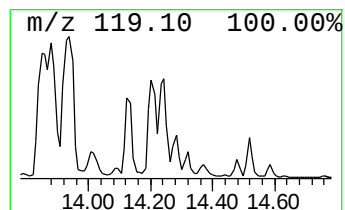
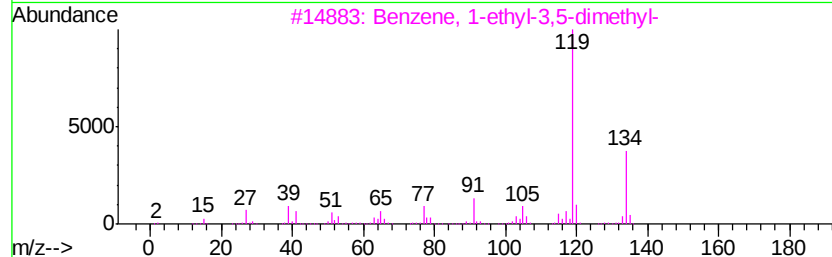
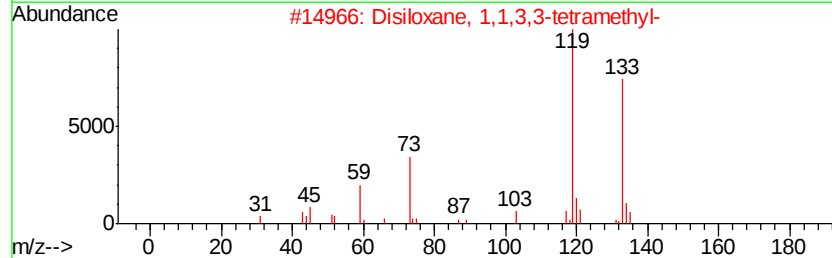
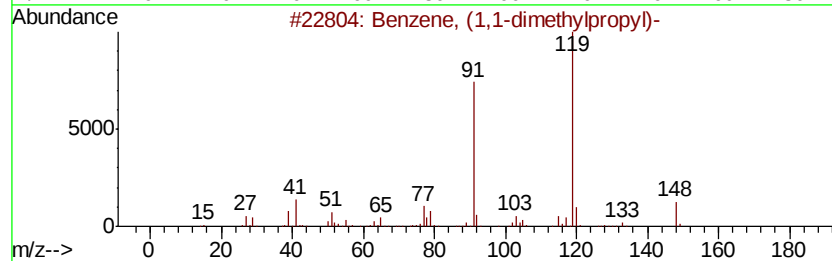
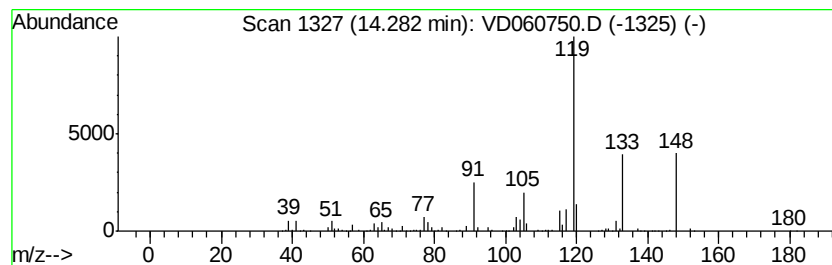
Instrument :
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ClientSampleId :
AOC2-06A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1,1-dimethylpropyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	147.92 ug/l	15631500	1,4-Dichlorobenzene-d4	13.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1,1-dimethylpropyl)-	148 C11H16	002049-95-8 50
2		Disiloxane, 1,1,3,3-tetramethyl-	134 C4H14OSi2	003277-26-7 47
3		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 47
4		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 47
5		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD122818\
Data File : VD060750.D
Acq On : 28 Dec 2018 16:59
Operator : VA/AP
Sample : J6590-01
Misc : 5.85g/5ml/MSVOA D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AOC2-06A

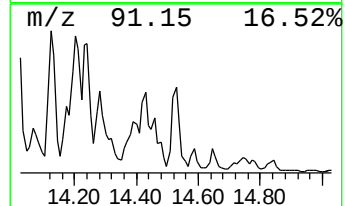
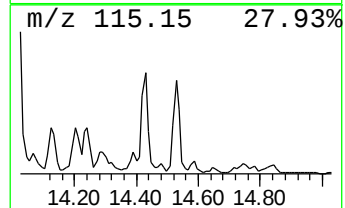
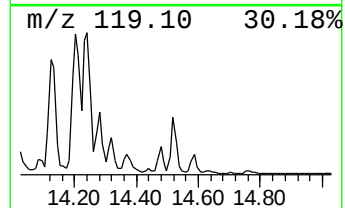
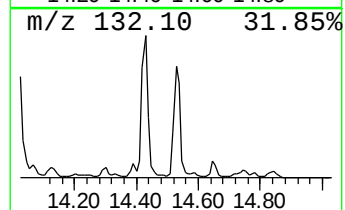
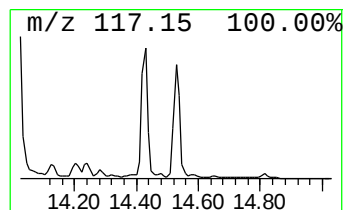
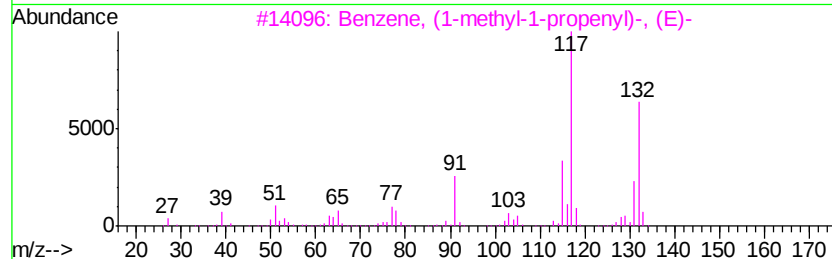
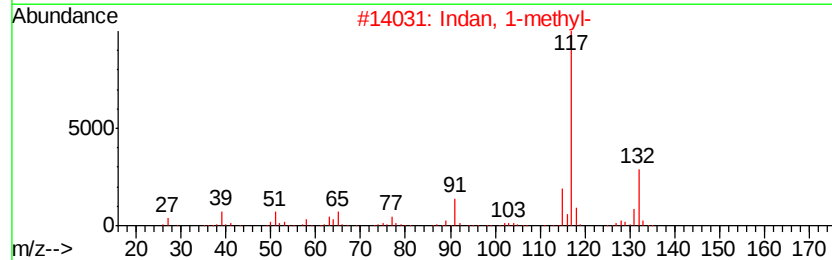
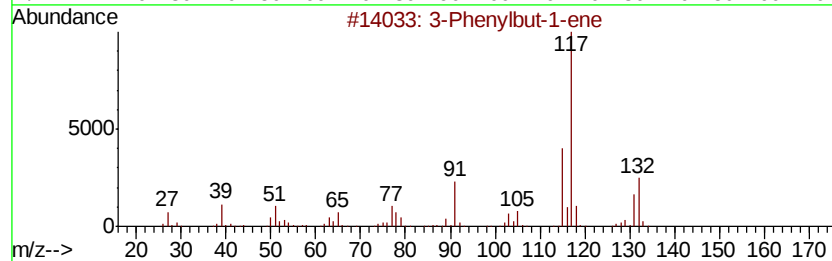
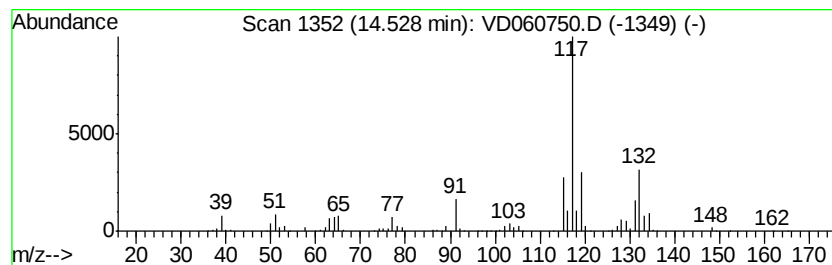
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Quant Title : SW846 8260

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

Peak Number 10 3-Phenylbut-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	159.23 ug/l	16826900	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
2		Indan, 1-methyl-	132	C10H12	000767-58-8	81
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD122818\
Data File : VD060750.D
Acq On : 28 Dec 2018 16:59
Operator : VA/AP
Sample : J6590-01
Misc : 5.85g/5ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AOC2-06A

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.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, 1,2,3-tr...	12.84	461.6	ug/l	48778300	4	13.40	5283910	50.0
Benzene, 1-ethyl-...	12.98	326.1	ug/l	34460900	4	13.40	5283910	50.0
Benzene, 1,4-diet...	13.59	216.7	ug/l	22897600	4	13.40	5283910	50.0
Benzene, (1-methy...	13.79	240.8	ug/l	25448300	4	13.40	5283910	50.0
Benzene, 2-ethyl-...	13.85	212.3	ug/l	22440900	4	13.40	5283910	50.0
Benzene, 1-ethyl-...	13.94	277.2	ug/l	29297700	4	13.40	5283910	50.0
Benzene, 1-methyl...	14.02	258.3	ug/l	27290900	4	13.40	5283910	50.0
Benzene, 1-ethyl-...	14.20	141.2	ug/l	14925200	4	13.40	5283910	50.0
Benzene, (1,1-dim...	14.28	147.9	ug/l	15631500	4	13.40	5283910	50.0
3-Phenylbut-1-ene	14.53	159.2	ug/l	16826900	4	13.40	5283910	50.0