

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF111918\
 Data File : VF060800.D
 Acq On : 20 Nov 2018 2:42
 Operator : VA/AP
 Sample : J5950-02
 Misc : 6.20g/5mL/MSVOA-F/SOIL
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 SSI-04

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_F\METHODS\82F111318S.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	1.081	41	46	53	rBV2	22403	58333	2.79%	0.130%
2	4.011	335	343	348	rBV2	12911	48772	2.34%	0.109%
3	4.120	348	354	366	rVB2	130896	450124	21.56%	1.005%
4	4.869	422	430	442	rBV2	257137	932169	44.66%	2.081%
5	5.461	482	490	498	rBV6	13040	54816	2.63%	0.122%
6	5.609	498	505	517	rVV	294115	797864	38.22%	1.781%
7	6.605	602	606	610	rBV3	10555	28354	1.36%	0.063%
8	7.039	644	650	658	rVB4	16990	64215	3.08%	0.143%
9	7.572	698	704	716	rVV	384733	939321	45.00%	2.097%
10	7.769	720	724	729	rVB3	10524	29405	1.41%	0.066%
11	8.292	773	777	783	rVB2	13406	31910	1.53%	0.071%
12	8.401	783	788	791	rBV5	13982	37657	1.80%	0.084%
13	8.470	791	795	803	rVB	44825	122456	5.87%	0.273%
14	8.637	805	812	817	rBV2	38181	105377	5.05%	0.235%
15	8.983	840	847	852	rBV4	28402	81190	3.89%	0.181%
16	9.131	857	862	867	rVB2	24643	69691	3.34%	0.156%
17	9.308	874	880	883	rBV	24459	56604	2.71%	0.126%
18	9.545	897	904	908	rBV4	11465	40690	1.95%	0.091%
19	9.653	911	915	917	rVV4	11098	29535	1.41%	0.066%
20	9.713	918	921	924	rVV5	17232	46175	2.21%	0.103%
21	9.791	924	929	938	rVV	391181	971396	46.53%	2.168%
22	9.910	938	941	946	rVV	37225	88634	4.25%	0.198%
23	10.245	969	975	983	rBV4	39768	136765	6.55%	0.305%
24	10.364	984	987	993	rVB5	12276	28863	1.38%	0.064%
25	10.531	995	1004	1011	rVB3	55741	153080	7.33%	0.342%
26	10.679	1015	1019	1028	rBV5	102608	282713	13.54%	0.631%
27	10.798	1028	1031	1033	rBV4	14373	29923	1.43%	0.067%
28	10.837	1033	1035	1038	rVB	20853	32879	1.58%	0.073%
29	10.965	1043	1048	1051	rVB5	13733	34149	1.64%	0.076%
30	11.064	1053	1058	1062	rBV5	34054	102076	4.89%	0.228%
31	11.133	1062	1065	1068	rBV	69970	128095	6.14%	0.286%
32	11.281	1076	1080	1084	rVB2	97722	190289	9.12%	0.425%
33	11.419	1088	1094	1100	rBV	398641	744764	35.68%	1.662%
34	11.498	1100	1102	1106	rVB4	21965	43611	2.09%	0.097%

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Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
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35	11.567	1106	1109	1111	rBV2	74318	128351	6.15%	0.286%
36	11.606	1111	1113	1118	rVB	146708	225007	10.78%	0.502%
37	11.735	1123	1126	1129	rBV2	59831	99481	4.77%	0.222%
38	11.833	1133	1136	1140	rVB	156486	232110	11.12%	0.518%
39	11.893	1140	1142	1145	rBV2	18130	32870	1.57%	0.073%
40	11.962	1145	1149	1151	rBV4	34351	75104	3.60%	0.168%
41	12.031	1151	1156	1159	rVB2	244311	514618	24.65%	1.149%
42	12.080	1159	1161	1166	rVB4	23521	45569	2.18%	0.102%
43	12.169	1166	1170	1171	rBV3	39654	80508	3.86%	0.180%
44	12.218	1171	1175	1179	rVB	997549	1577075	75.55%	3.520%
45	12.317	1180	1185	1190	rVB	166836	287182	13.76%	0.641%
46	12.396	1190	1193	1195	rBV2	57562	101768	4.88%	0.227%
47	12.435	1195	1197	1198	rBV	30009	37224	1.78%	0.083%
48	12.465	1198	1200	1206	rVB	138590	236415	11.33%	0.528%
49	12.563	1206	1210	1213	rBV	564408	916895	43.92%	2.047%
50	12.623	1213	1216	1220	rVB	472426	724611	34.71%	1.617%
51	12.731	1224	1227	1231	rBV3	236638	527941	25.29%	1.178%
52	12.820	1231	1236	1237	rBV	210389	409081	19.60%	0.913%
53	12.849	1237	1239	1243	rVB2	290785	494405	23.68%	1.104%
54	12.938	1245	1248	1253	rBV4	99178	189989	9.10%	0.424%
55	13.017	1253	1256	1260	rVB	349043	517038	24.77%	1.154%
56	13.086	1260	1263	1268	rBV2	485285	1269116	60.80%	2.833%
57	13.165	1268	1271	1276	rBV2	558853	874938	41.91%	1.953%
58	13.234	1276	1278	1283	rVB2	276641	561183	26.88%	1.253%
59	13.313	1283	1286	1288	rBV3	121420	258766	12.40%	0.578%
60	13.343	1288	1289	1292	rVB2	80475	100143	4.80%	0.224%
61	13.392	1292	1294	1296	rVB	137883	165982	7.95%	0.370%
62	13.431	1296	1298	1303	rBV	216392	302235	14.48%	0.675%
63	13.510	1303	1306	1308	rBV	346979	516019	24.72%	1.152%
64	13.550	1308	1310	1312	rBV	526105	733927	35.16%	1.638%
65	13.629	1316	1318	1319	rBV	61099	86669	4.15%	0.193%
66	13.648	1319	1320	1324	rVB2	141657	142433	6.82%	0.318%
67	13.717	1324	1327	1330	rBV	650670	966711	46.31%	2.158%
68	13.767	1330	1332	1335	rVB3	190564	301730	14.45%	0.673%
69	13.826	1335	1338	1339	rBV	270643	398523	19.09%	0.890%
70	13.865	1339	1342	1346	rBV3	948110	1749415	83.81%	3.905%
71	13.925	1346	1348	1351	rVB2	343927	496672	23.79%	1.109%

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72	13.984	1351	1354	1357	rVB2	551245	842560	40.36%	1.881%
73	14.063	1359	1362	1363	rBV	249179	396938	19.02%	0.886%
74	14.092	1363	1365	1367	rVB	382293	500267	23.97%	1.117%
75	14.161	1367	1372	1375	rBV4	639105	1472395	70.54%	3.286%
76	14.230	1377	1379	1383	rVB2	443564	622170	29.81%	1.389%
77	14.309	1385	1387	1389	rBV2	211709	244281	11.70%	0.545%
78	14.368	1389	1393	1395	rBV3	138551	233003	11.16%	0.520%
79	14.418	1395	1398	1401	rBV3	495641	819175	39.24%	1.828%
80	14.467	1401	1403	1406	rBV	489180	706373	33.84%	1.577%
81	14.576	1411	1414	1415	rVV2	221238	412149	19.74%	0.920%
82	14.605	1415	1417	1424	rVV2	448574	1008494	48.31%	2.251%
83	14.714	1426	1428	1429	rVV	147251	247666	11.86%	0.553%
84	14.753	1429	1432	1435	rVV	1220718	2087460	100.00%	4.659%
85	14.832	1438	1440	1441	rVV	215043	303163	14.52%	0.677%
86	14.881	1441	1445	1450	rVV3	362895	1278436	61.24%	2.854%
87	14.950	1450	1452	1455	rVV2	144054	244913	11.73%	0.547%
88	15.010	1455	1458	1463	rVV2	708532	1500917	71.90%	3.350%
89	15.118	1466	1469	1472	rVV	793288	1407793	67.44%	3.142%
90	15.177	1472	1475	1481	rVV4	346320	954922	45.75%	2.131%
91	15.296	1483	1487	1490	rVV	888515	1675932	80.29%	3.741%
92	15.355	1490	1493	1496	rVV2	379331	933576	44.72%	2.084%
93	15.414	1496	1499	1503	rVV	545613	1003379	48.07%	2.240%
94	15.503	1506	1508	1510	rVV2	75346	143110	6.86%	0.319%
95	15.572	1513	1515	1517	rVV	140837	214618	10.28%	0.479%
96	15.651	1520	1523	1528	rVV	394207	797062	38.18%	1.779%
97	15.720	1528	1530	1536	rVV5	98766	239434	11.47%	0.534%
98	15.828	1536	1541	1545	rVB3	104253	307575	14.73%	0.687%
99	15.947	1550	1553	1555	rBV2	97516	199249	9.55%	0.445%
100	16.006	1555	1559	1567	rVB4	236579	635559	30.45%	1.419%

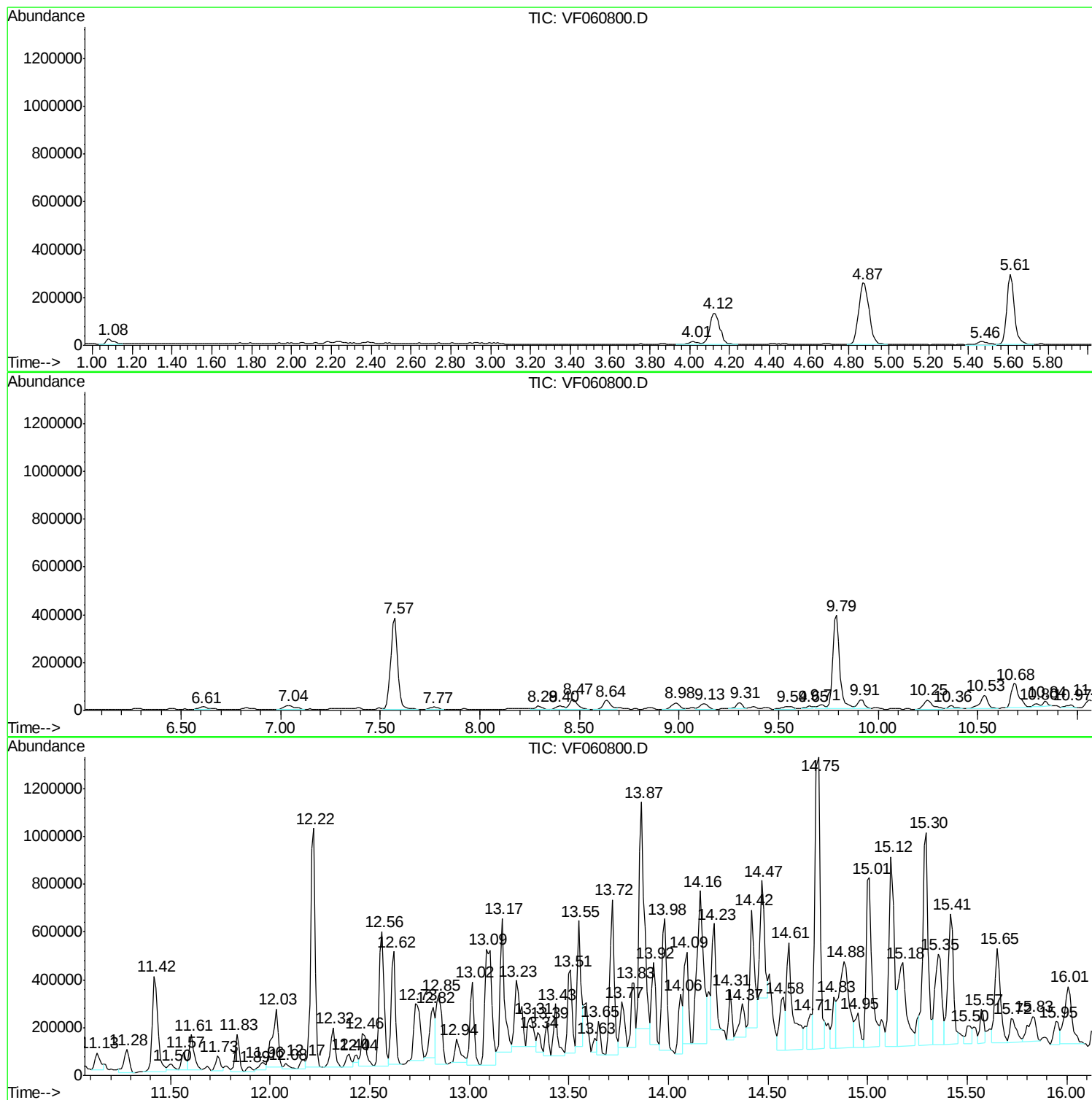
Sum of corrected areas: 44802168

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



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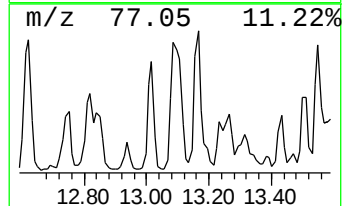
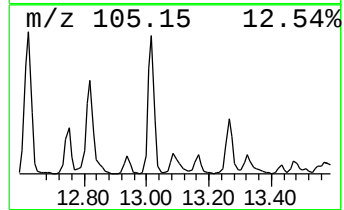
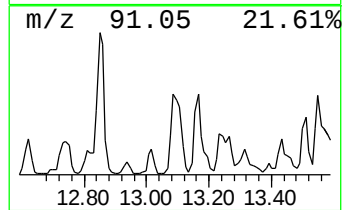
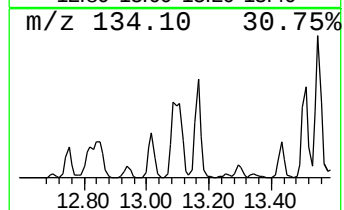
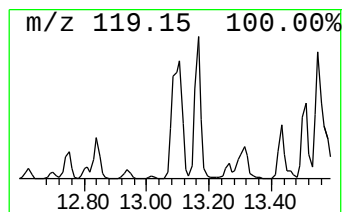
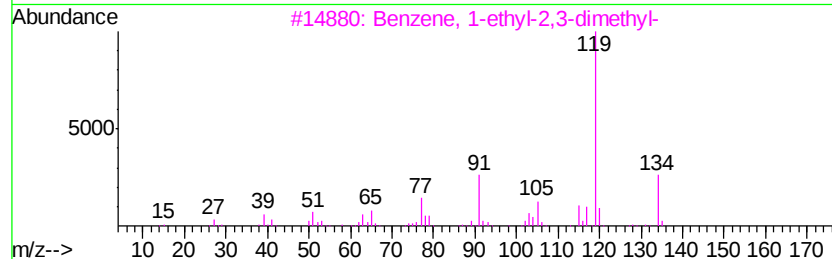
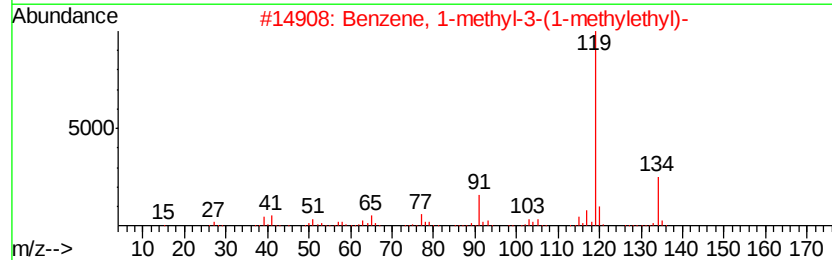
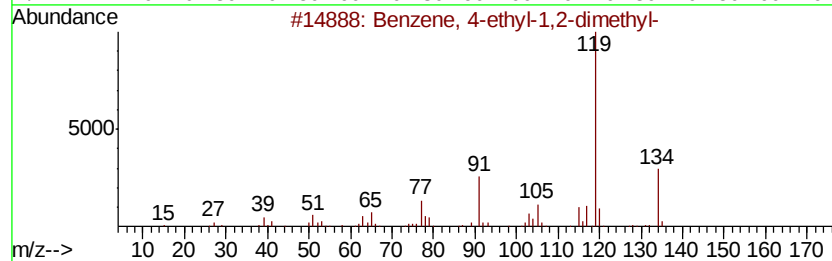
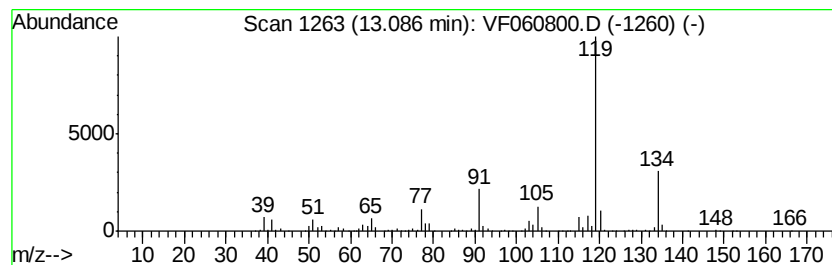
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.09	69.21 ug/l	1269120	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		p-Cymene	134	C10H14	000099-87-6	94
5		o-Cymene	134	C10H14	000527-84-4	94



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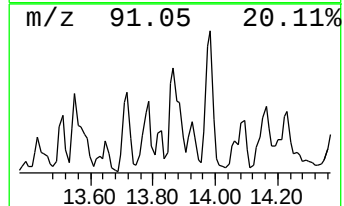
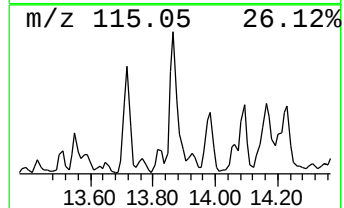
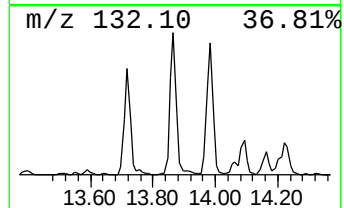
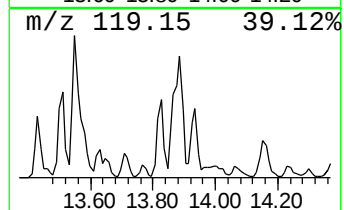
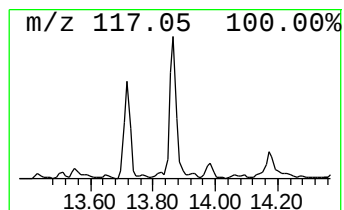
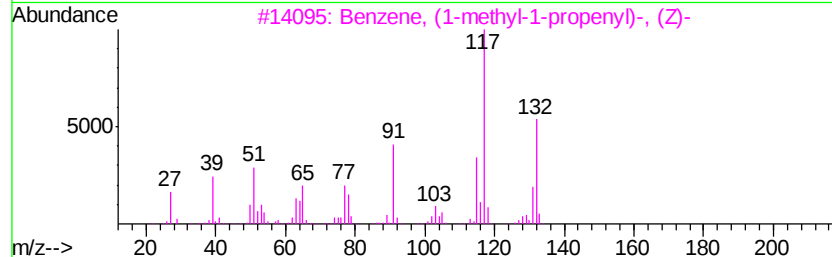
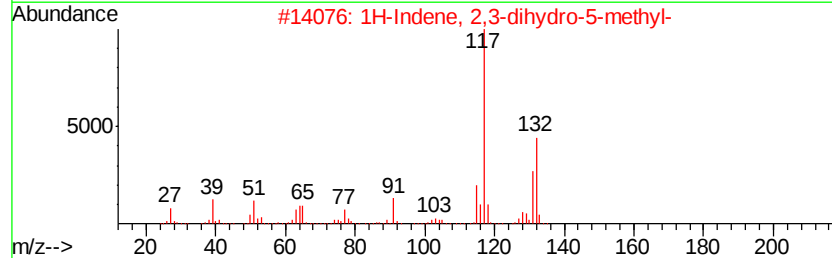
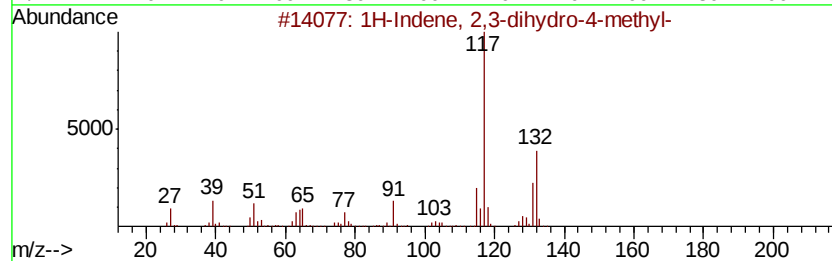
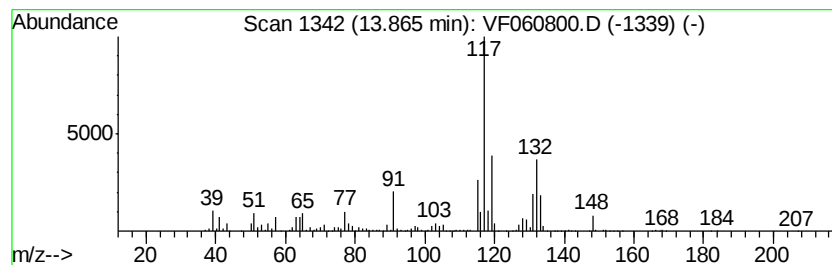
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Peak Number 2 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	95.40 ug/l	1749420	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	68
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64



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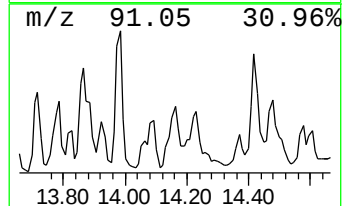
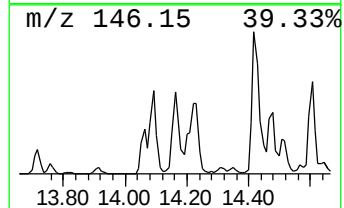
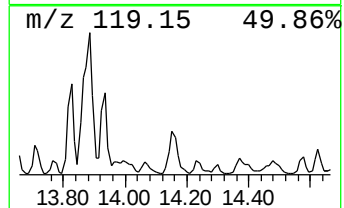
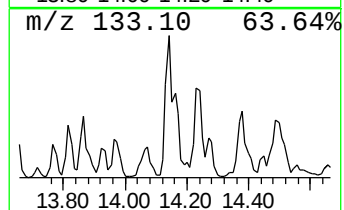
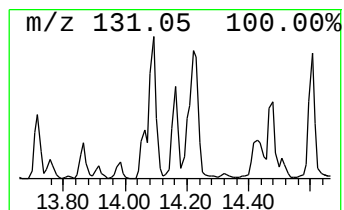
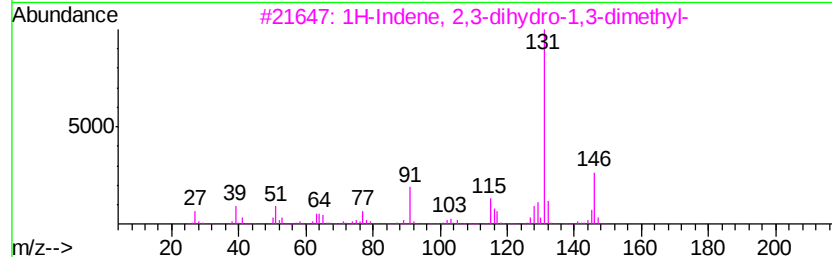
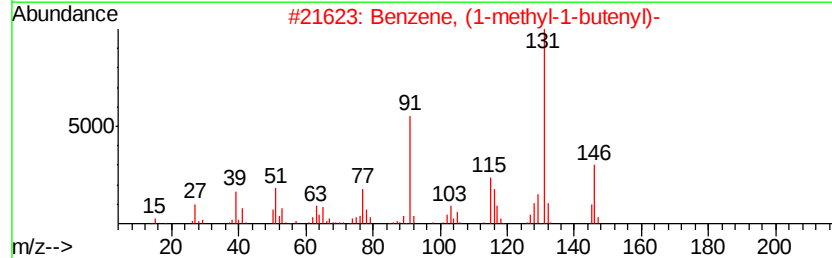
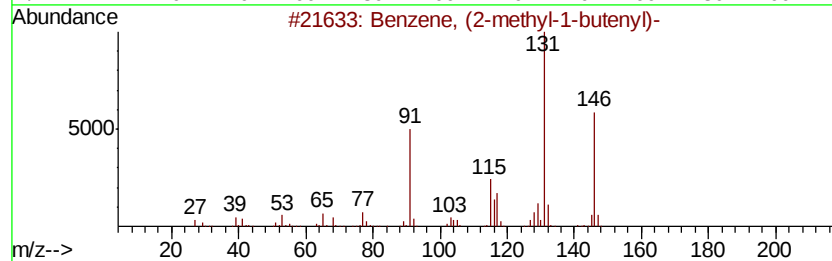
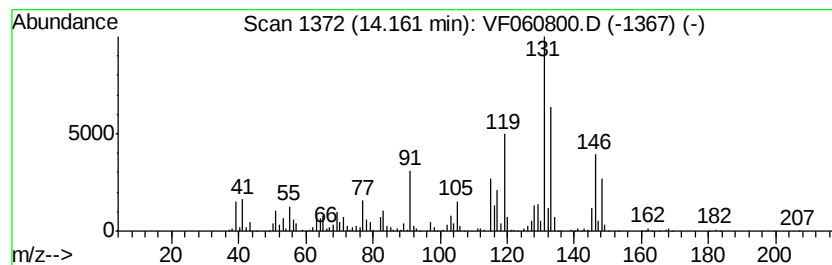
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.16	80.29 ug/l	1472400	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	92
2		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	89
3		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060800.D
Acq On : 20 Nov 2018 2:42
Operator : VA/AP
Sample : J5950-02
Misc : 6.20µ/5mL/MSVOA-F/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-04

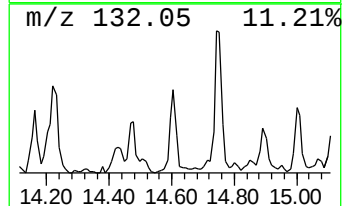
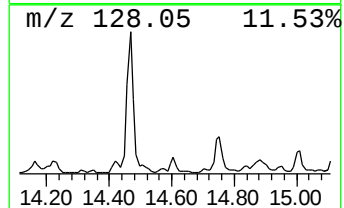
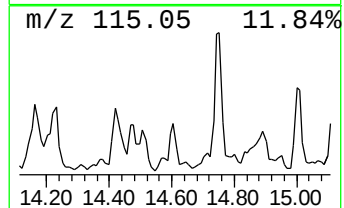
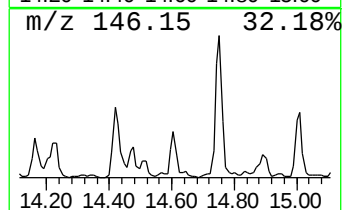
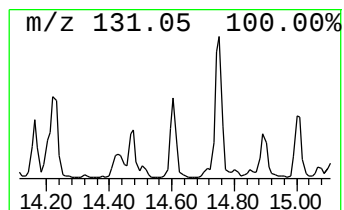
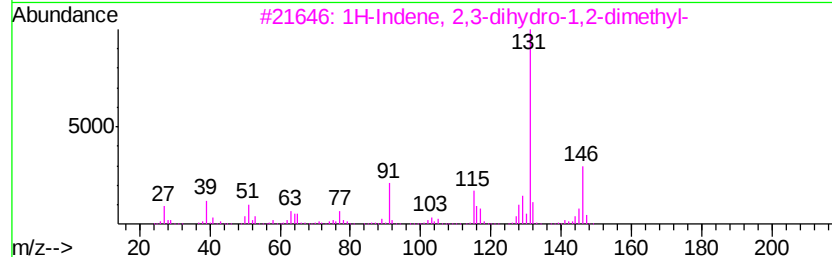
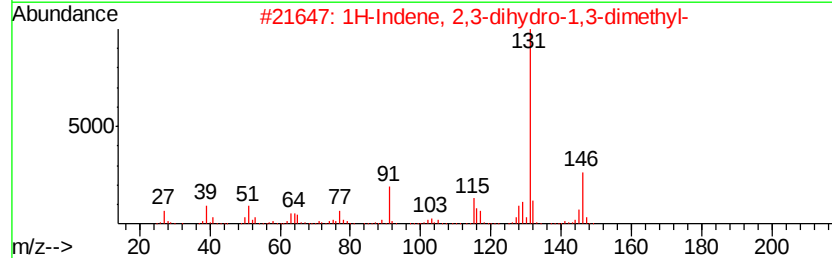
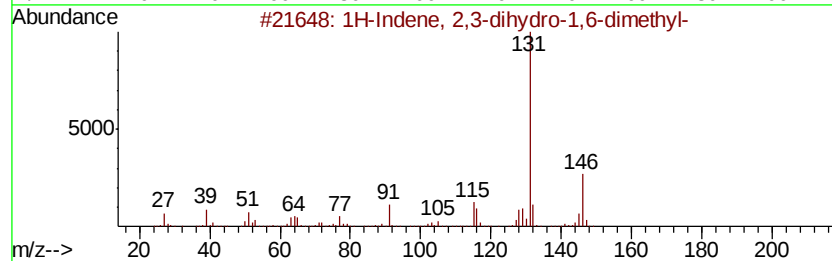
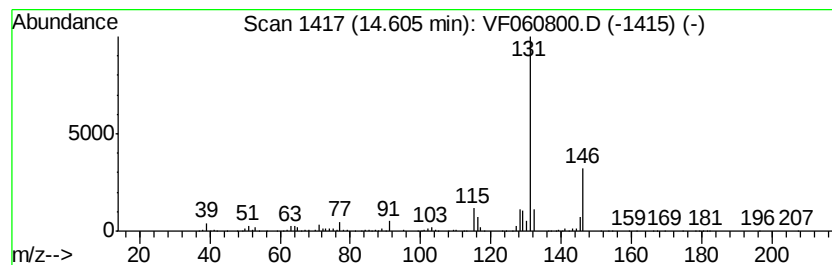
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Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	55.00 ug/l	1008490	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	91
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060800.D
Acq On : 20 Nov 2018 2:42
Operator : VA/AP
Sample : J5950-02
Misc : 6.20g/5mL/MSVOA-F/SOIL
ALS Vial : 24 Sample Multiplier: 1

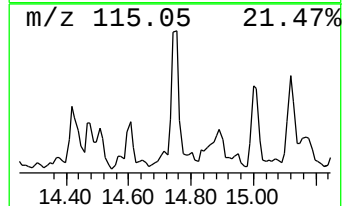
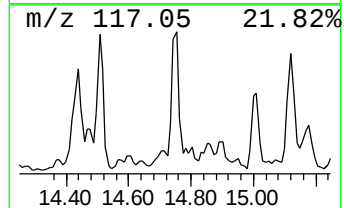
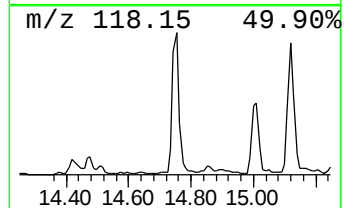
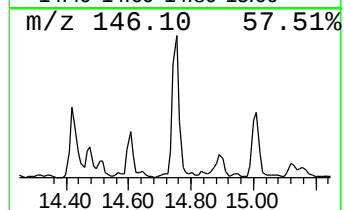
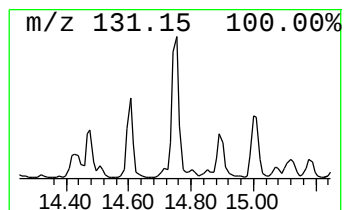
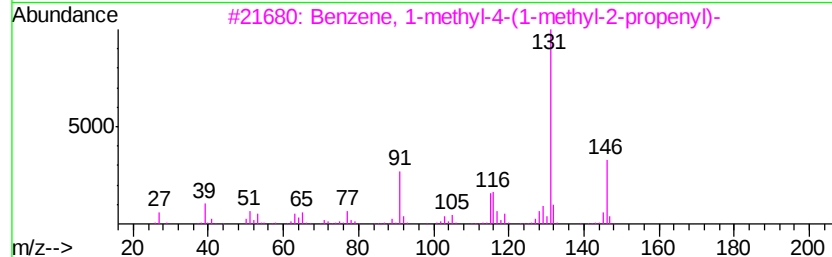
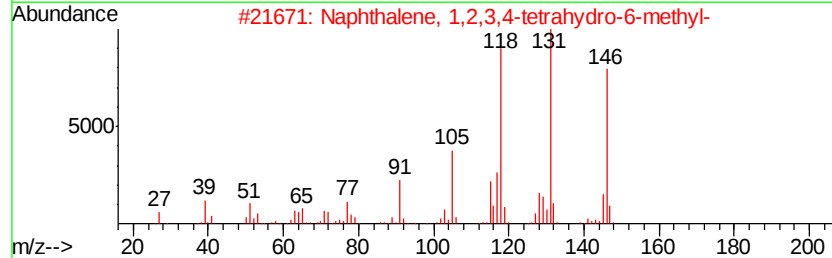
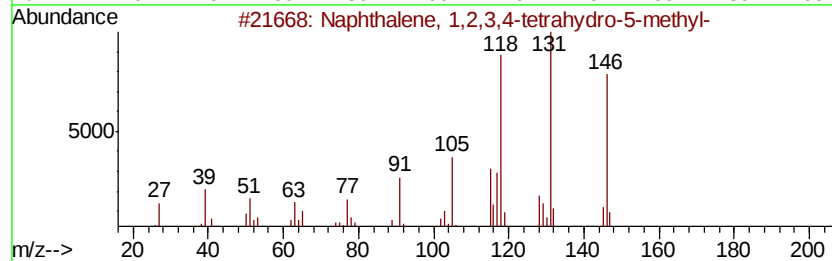
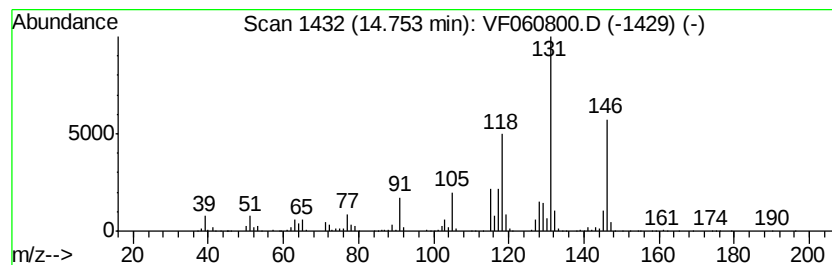
Instrument :
MSVOA_F
ClientSampleId :
SSI-04

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetra... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	113.83 ug/l	2087460	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 96
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 94
3		Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1 70
4		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 68
5		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060800.D
Acq On : 20 Nov 2018 2:42
Operator : VA/AP
Sample : J5950-02
Misc : 6.20g/5mL/MSVOA-F/SOIL
ALS Vial : 24 Sample Multiplier: 1

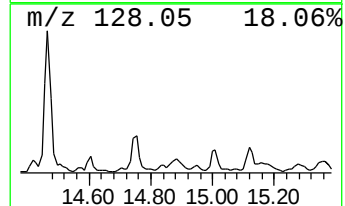
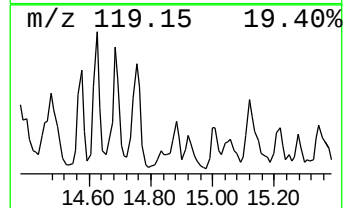
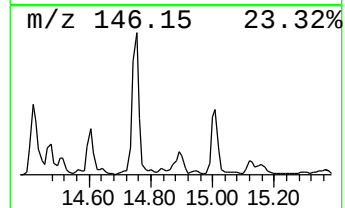
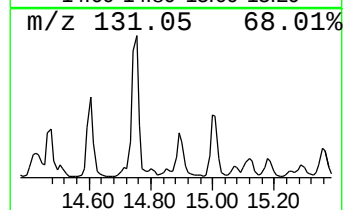
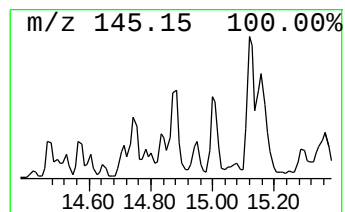
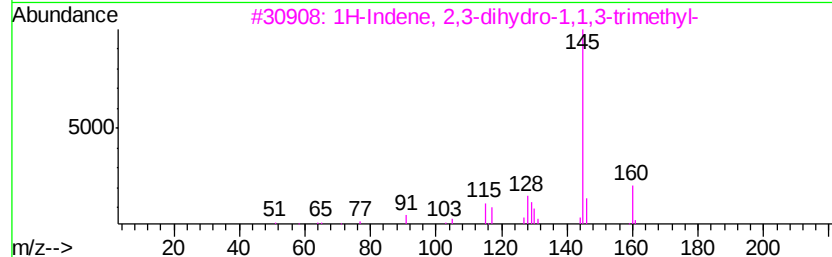
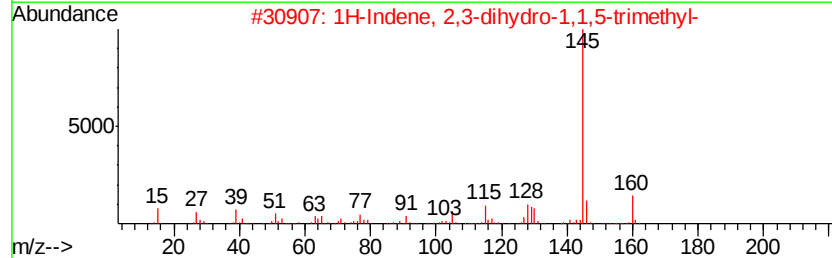
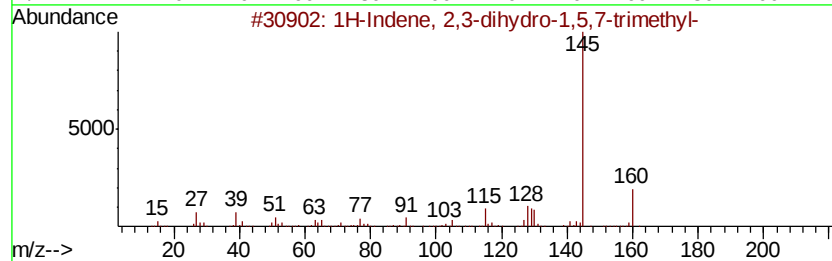
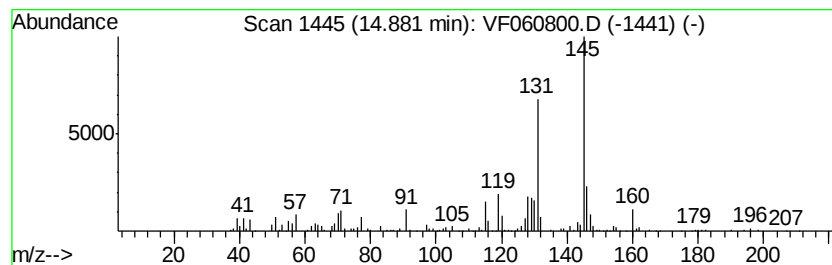
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MSVOA_F
ClientSampleId :
SSI-04

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-1,5,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	69.72 ug/l	1278440	1,4-Dichlorobenzene-d4	12.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,5,7-tri...	160 C12H16	054340-88-4	53
2	1H-Indene, 2,3-dihydro-1,1,5-tri...	160 C12H16	040650-41-7	53
3	1H-Indene, 2,3-dihydro-1,1,3-tri...	160 C12H16	002613-76-5	52
4	Benzene, 4-(2-butenyl)-1,2-dimet...	160 C12H16	054340-86-2	52
5	Benzene, 1-(1-methylethenyl)-3-(...	160 C12H16	001129-29-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060800.D
Acq On : 20 Nov 2018 2:42
Operator : VA/AP
Sample : J5950-02
Misc : 6.20g/5mL/MSVOA-F/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-04

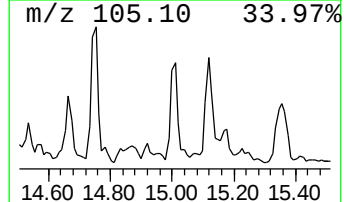
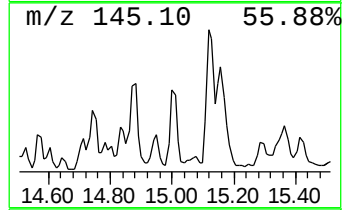
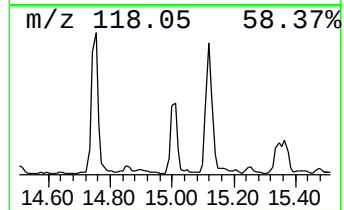
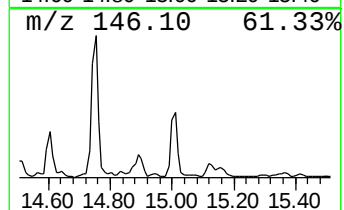
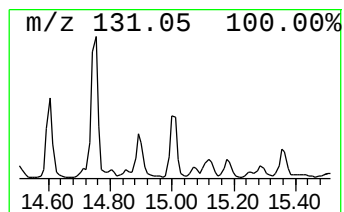
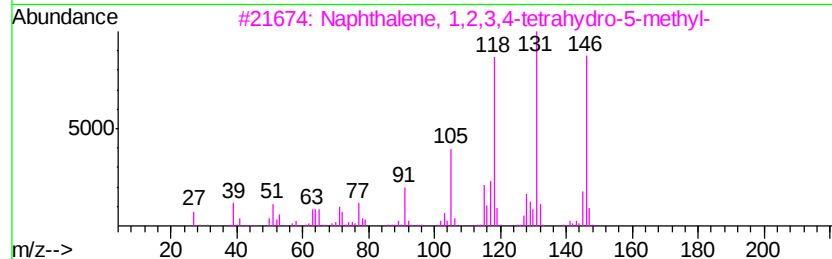
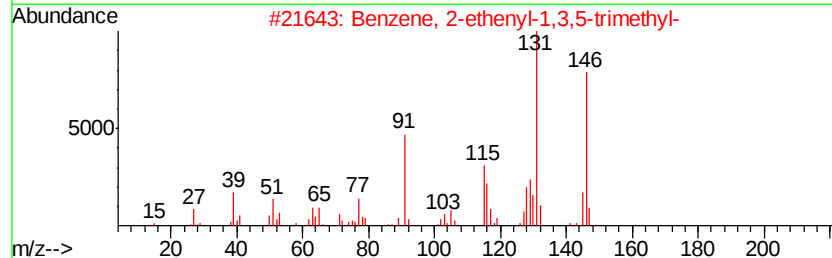
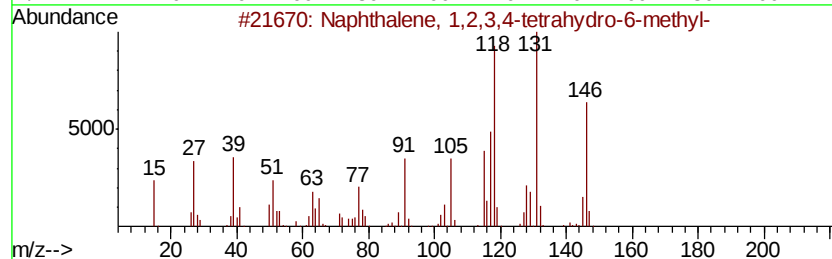
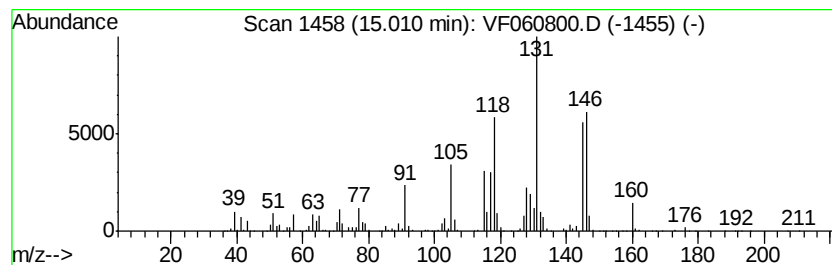
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.01	81.85 ug/l	1500920	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
2		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	83
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	76
4		1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	70
5		2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060800.D
Acq On : 20 Nov 2018 2:42
Operator : VA/AP
Sample : J5950-02
Misc : 6.20g/5mL/MSVOA-F/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-04

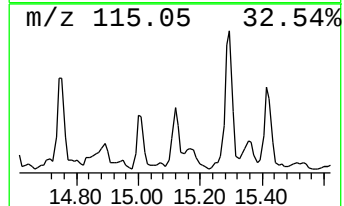
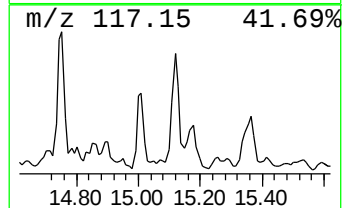
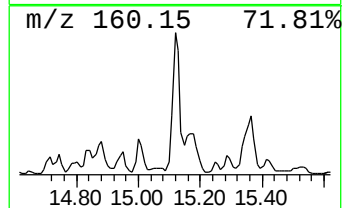
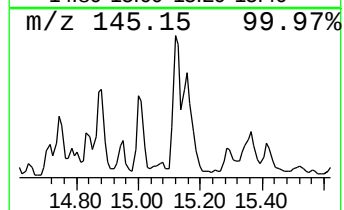
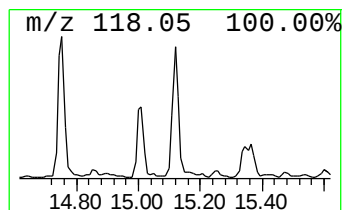
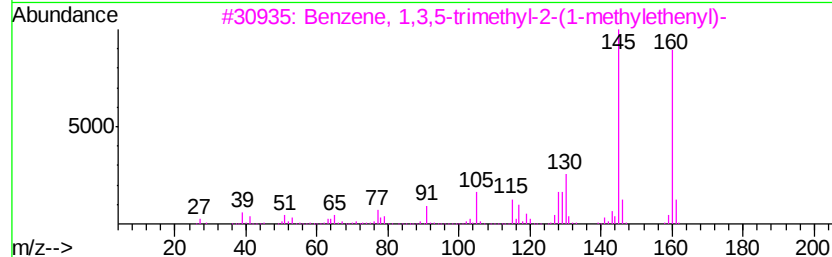
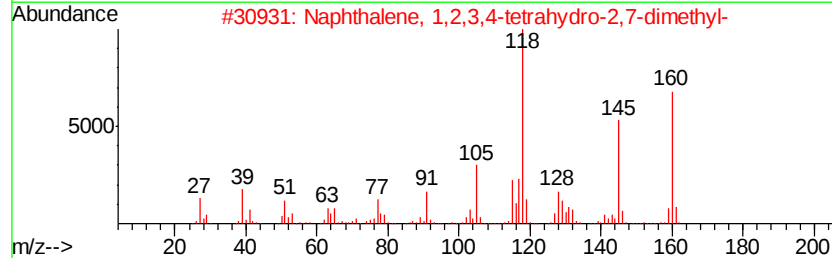
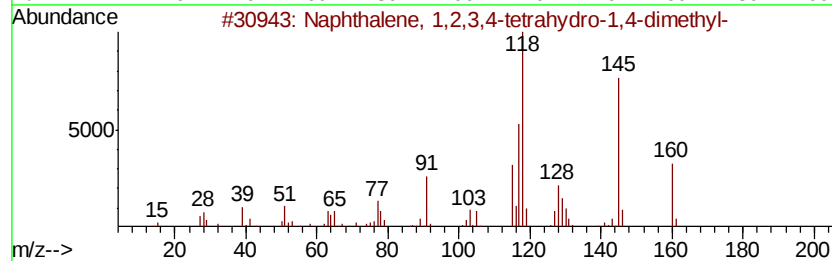
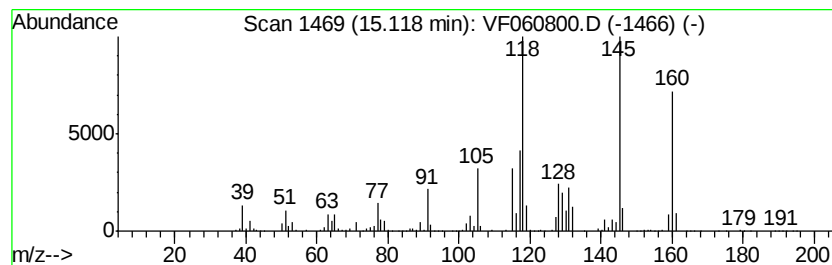
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	76.77 ug/l	1407790	1,4-Dichlorobenzene-d4	12.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	97
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	95
3		Benzene, 1,3,5-trimethyl-2-(1-methyl-...	160	C12H16	014679-13-1	78
4		Benzene, 1,2,4-trimethyl-5-(1-methyl-...	160	C12H16	054340-84-0	70
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF111918\
Data File : VF060800.D
Acq On : 20 Nov 2018 2:42
Operator : VA/AP
Sample : J5950-02
Misc : 6.20g/5mL/MSVOA-F/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SSI-04

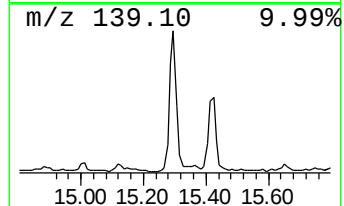
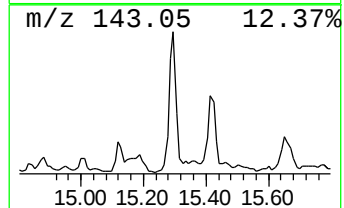
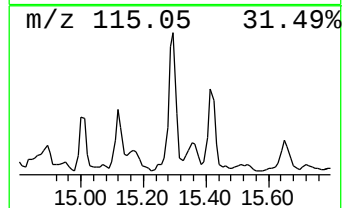
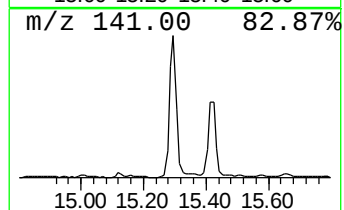
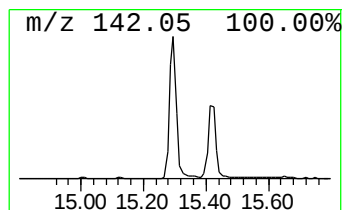
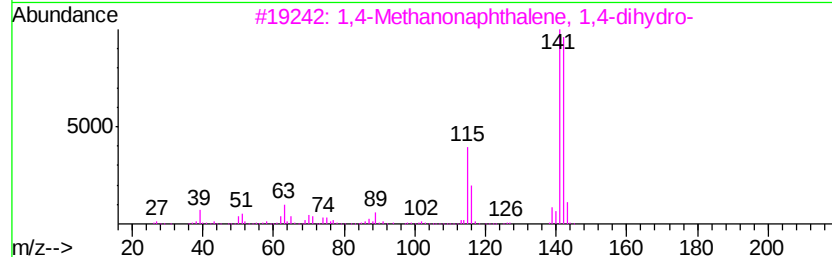
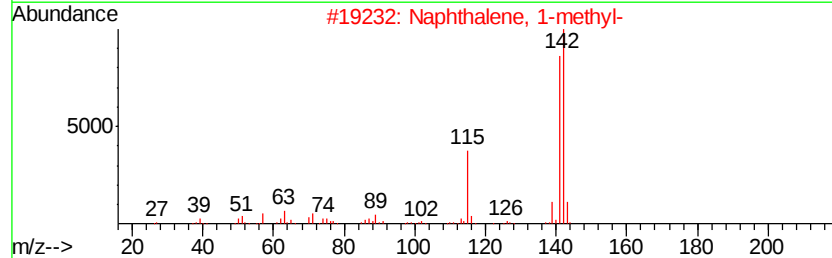
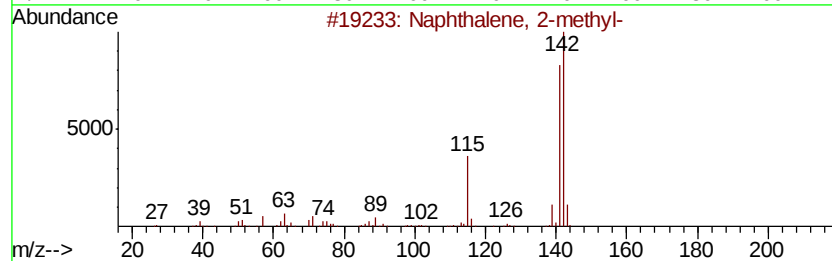
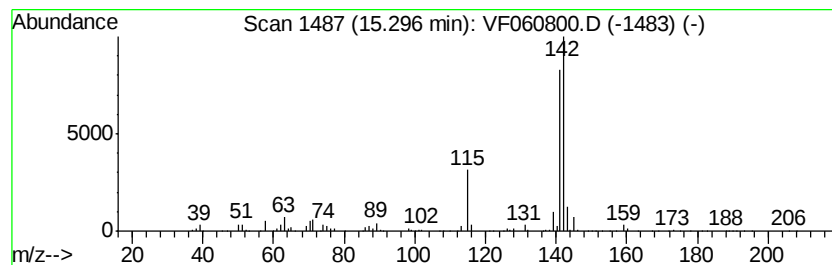
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.30	91.39 ug/l	1675930	1,4-Dichlorobenzene-d4	12.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
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	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF111918\
Data File : VF060800.D
Acq On : 20 Nov 2018 2:42
Operator : VA/AP
Sample : J5950-02
Misc : 6.20g/5mL/MSVOA-F/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
SSI-04

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.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, 4-ethyl-...	13.09	69.2	ug/l	1269120	4	12.56	916895	50.0
1H-Indene, 2,3-di...	13.87	95.4	ug/l	1749420	4	12.56	916895	50.0
Benzene, (2-methy...	14.16	80.3	ug/l	1472400	4	12.56	916895	50.0
1H-Indene, 2,3-di...	14.61	55.0	ug/l	1008490	4	12.56	916895	50.0
Naphthalene, 1,2,...	14.75	113.8	ug/l	2087460	4	12.56	916895	50.0
1H-Indene, 2,3-di...	14.88	69.7	ug/l	1278440	4	12.56	916895	50.0
Naphthalene, 1,2,...	15.01	81.8	ug/l	1500920	4	12.56	916895	50.0
Naphthalene, 1,2,...	15.12	76.8	ug/l	1407790	4	12.56	916895	50.0
Naphthalene, 1-me...	15.30	91.4	ug/l	1675930	4	12.56	916895	50.0