

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD062218\
 Data File : VD058309.D
 Acq On : 22 Jun 2018 21:12
 Operator : VA/AP
 Sample : J3663-12
 Misc : 8.94g/5ml/MSVOA D/SOIL
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 WP021SB470-27-29-180618

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\82D061918S.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	1.213	93	100	102	rBV	1458725	2929873	4.81%	0.378%
2	1.272	102	106	127	rVB	1974649	6294244	10.34%	0.813%
3	5.544	529	540	545	rBV	423495	1267406	2.08%	0.164%
4	5.643	545	550	553	rVV	634335	2001991	3.29%	0.259%
5	5.692	553	555	567	rVB	382757	1187193	1.95%	0.153%
6	6.017	580	588	595	rBV2	360951	1000401	1.64%	0.129%
7	6.214	595	608	633	rVB	3953778	18496144	30.39%	2.388%
8	6.686	650	656	667	rBV	844086	2273616	3.74%	0.294%
9	6.962	677	684	692	rBV	616320	2121385	3.49%	0.274%
10	7.336	705	722	726	rBV5	1493557	8078045	13.27%	1.043%
11	7.405	726	729	734	rVV2	1279506	3891269	6.39%	0.502%
12	7.484	734	737	748	rVB	564274	1945858	3.20%	0.251%
13	7.671	748	756	771	rVB3	597043	2698976	4.43%	0.349%
14	7.907	771	780	790	rBV	3447240	13001067	21.36%	1.679%
15	8.084	790	798	818	rVB	3987307	17817319	29.27%	2.301%
16	8.380	818	828	835	rBV2	679181	2969265	4.88%	0.383%
17	8.576	842	848	850	rBV3	178166	617784	1.02%	0.080%
18	8.665	850	857	865	rBV2	2907174	14613784	24.01%	1.887%
19	8.941	876	885	898	rVB3	1407742	6529047	10.73%	0.843%
20	9.216	906	913	923	rBV	3448509	12090022	19.86%	1.561%
21	9.374	923	929	934	rBV	1552833	4761751	7.82%	0.615%
22	9.462	934	938	940	rBV	928947	2268599	3.73%	0.293%
23	9.521	940	944	949	rVV	1737416	5528688	9.08%	0.714%
24	9.600	949	952	958	rVB	1248947	3750430	6.16%	0.484%
25	9.758	958	968	978	rVB	5696635	23495636	38.60%	3.034%
26	9.925	978	985	1001	rBV2	6701969	39475363	64.86%	5.097%
27	10.132	1001	1006	1010	rBV	3891911	12084511	19.85%	1.560%
28	10.211	1010	1014	1020	rVB3	2817421	8863316	14.56%	1.145%
29	10.309	1021	1024	1027	rBV	416053	937892	1.54%	0.121%
30	10.447	1032	1038	1049	rVB	7642900	27066236	44.47%	3.495%
31	10.693	1054	1063	1068	rBV4	2380973	10351272	17.01%	1.337%
32	10.791	1068	1073	1079	rVB3	2818451	10964502	18.01%	1.416%
33	10.900	1079	1084	1089	rBV	4242723	14133433	23.22%	1.825%
34	11.008	1089	1095	1097	rBV2	3872643	12028613	19.76%	1.553%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	11.067	1097	1101	1104	rBV2	6899067	18616855	30.59%	2.404%
36	11.116	1104	1106	1113	rVB	9613433	28564953	46.93%	3.689%
37	11.225	1113	1117	1122	rBV	3411503	8245771	13.55%	1.065%
38	11.313	1122	1126	1128	rBV2	3632908	9806874	16.11%	1.266%
39	11.402	1128	1135	1140	rBV6	11530786	57509418	94.49%	7.426%
40	11.589	1147	1154	1159	rBV6	9784403	51202616	84.13%	6.612%
41	11.707	1163	1166	1180	rVB6	11250054	60864197	100.00%	7.859%
42	11.874	1180	1183	1184	rBV2	6506635	12696293	20.86%	1.639%
43	12.012	1194	1197	1200	rBV3	4446566	12255128	20.14%	1.583%
44	12.248	1216	1221	1224	rBV6	5806459	19645800	32.28%	2.537%
45	12.563	1248	1253	1256	rBV5	5346648	18433560	30.29%	2.380%
46	12.790	1271	1276	1279	rBV3	8310380	23556681	38.70%	3.042%
47	12.839	1279	1281	1283	rVB	4463942	6018207	9.89%	0.777%
48	13.253	1317	1323	1327	rBV6	9009933	25750924	42.31%	3.325%
49	13.508	1348	1349	1352	rVB	4313367	5227030	8.59%	0.675%
50	13.568	1352	1355	1358	rVV4	11663002	25742318	42.29%	3.324%
51	13.617	1358	1360	1361	rVV	3706202	5034892	8.27%	0.650%
52	13.646	1361	1363	1364	rVV	4772380	6933083	11.39%	0.895%
53	13.676	1364	1366	1367	rVV	8523660	12399944	20.37%	1.601%
54	13.715	1367	1370	1377	rVV5	12763311	41658525	68.45%	5.379%
55	13.804	1377	1379	1381	rVV	5604412	7793771	12.81%	1.006%
56	13.833	1381	1382	1385	rVV2	6140186	7919842	13.01%	1.023%
57	13.873	1385	1386	1390	rVB3	2729556	4248804	6.98%	0.549%
58	13.932	1390	1392	1394	rVB2	1069655	1494264	2.46%	0.193%
59	13.981	1394	1397	1399	rVB3	366634	609865	1.00%	0.079%
60	14.070	1404	1406	1408	rVB2	1210625	1447661	2.38%	0.187%
61	14.217	1418	1421	1423	rVB	1849631	2550101	4.19%	0.329%
62	14.276	1425	1427	1430	rVB	482463	645350	1.06%	0.083%

Sum of corrected areas: 774407658

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



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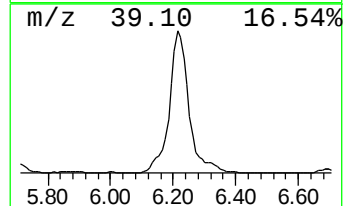
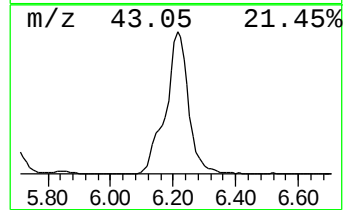
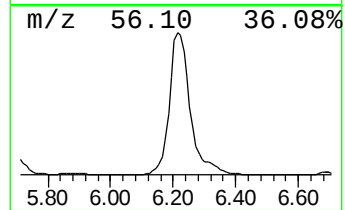
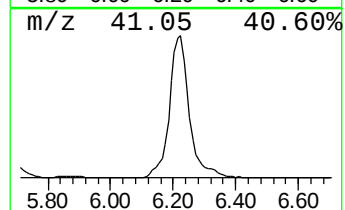
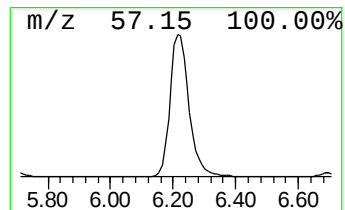
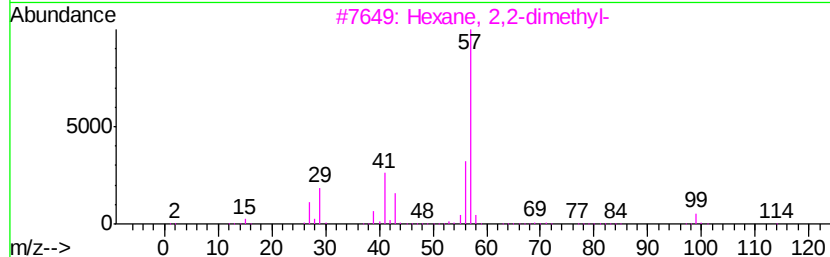
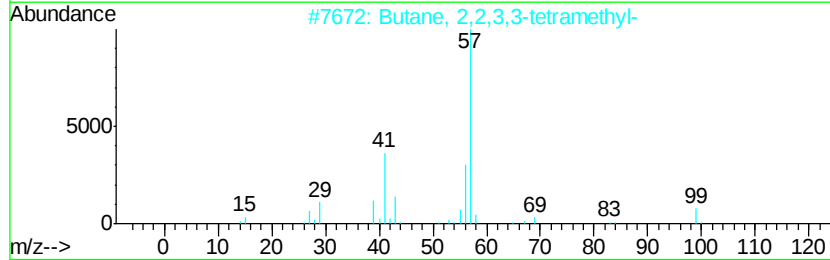
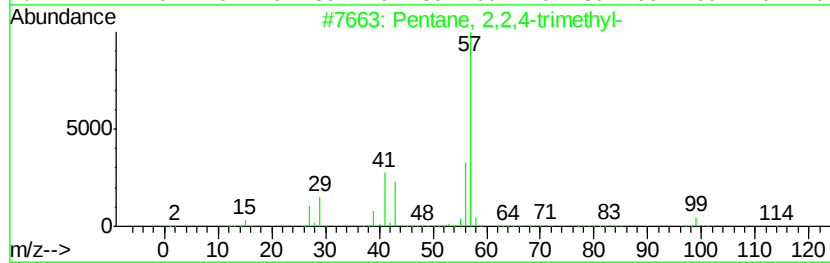
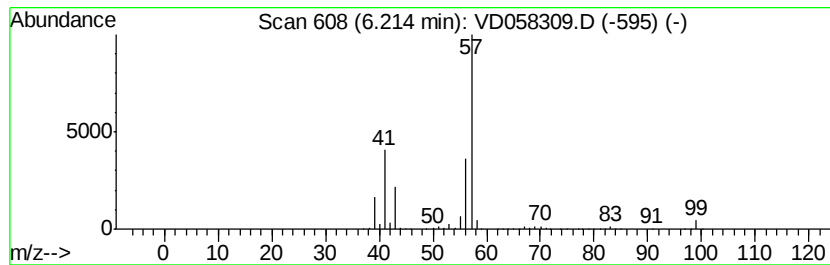
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.21	406.76 ug/l	18496100	1,4-Difluorobenzene	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	64
3		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
4		Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	42
5		Decane, 2,2,6-trimethyl-	184	C13H28	062237-97-2	42



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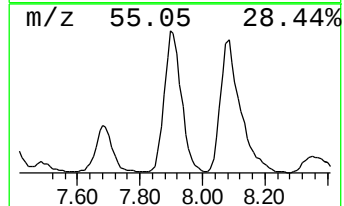
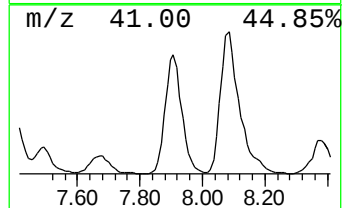
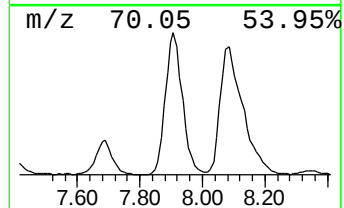
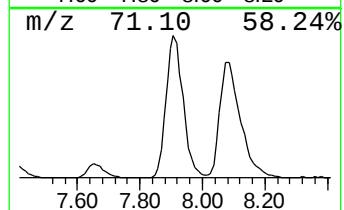
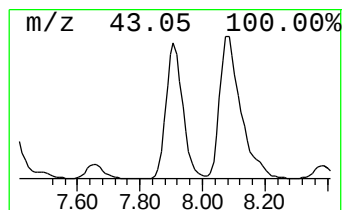
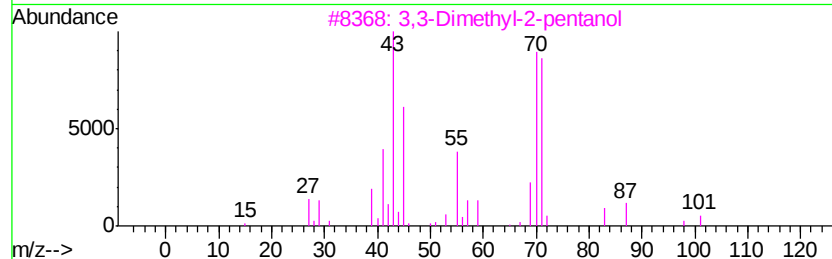
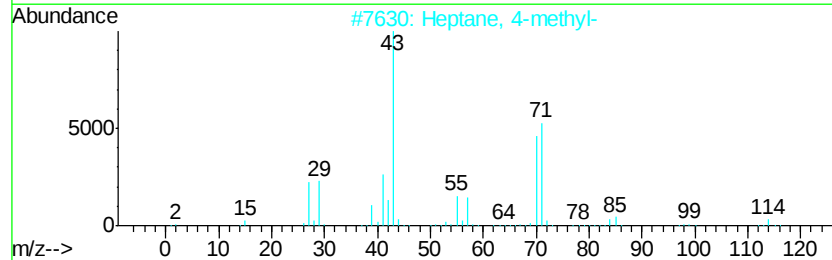
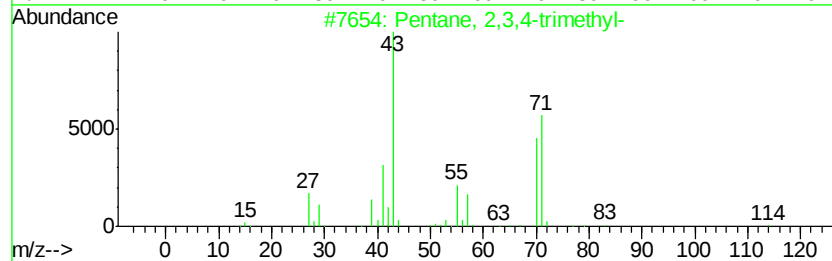
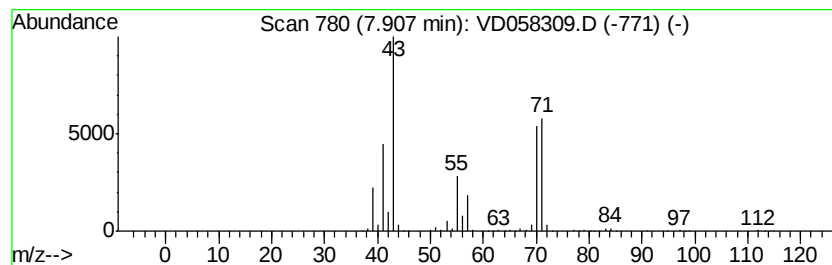
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2,3,4-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.91	285.91 ug/l	13001100	1,4-Difluorobenzene	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3		3,3-Dimethyl-2-pentanol	116	C7H16O	019781-24-9	64
4		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	64
5		Pentane, 3-ethyl-	100	C7H16	000617-78-7	64



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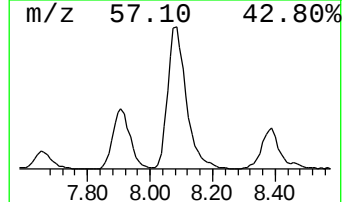
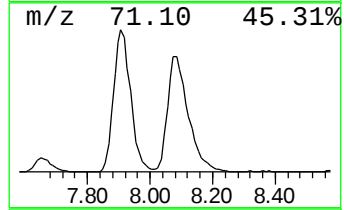
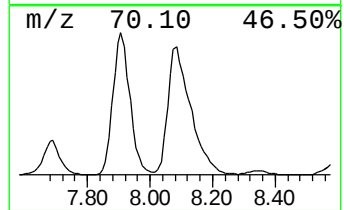
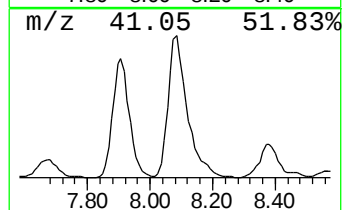
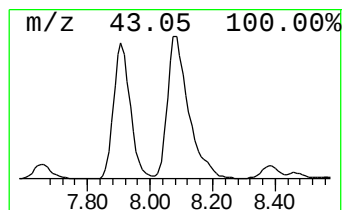
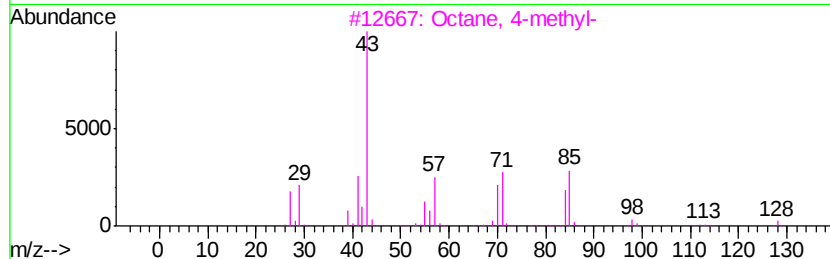
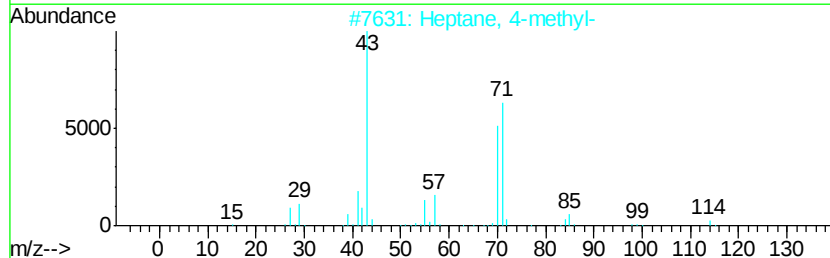
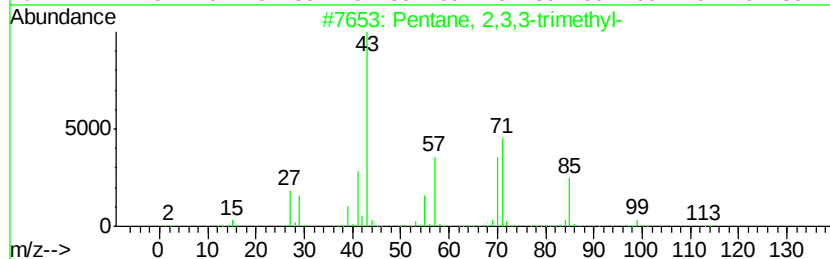
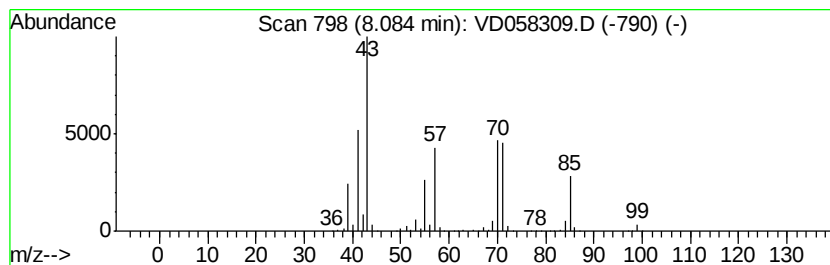
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.08	391.83 ug/l	17817300	1,4-Difluorobenzene	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3		Octane, 4-methyl-	128	C9H20	002216-34-4	64
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	64
5		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59



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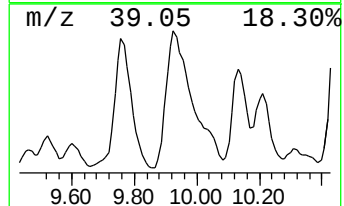
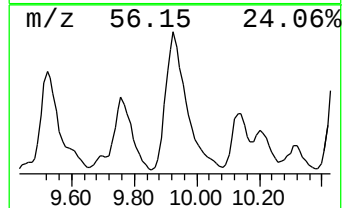
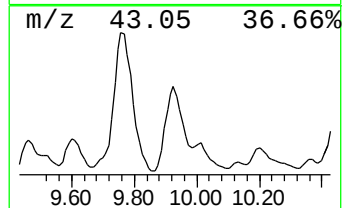
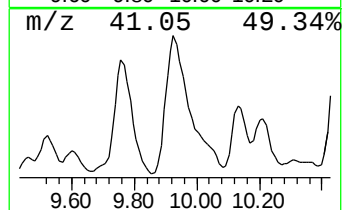
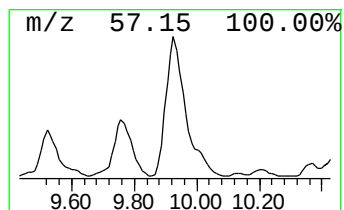
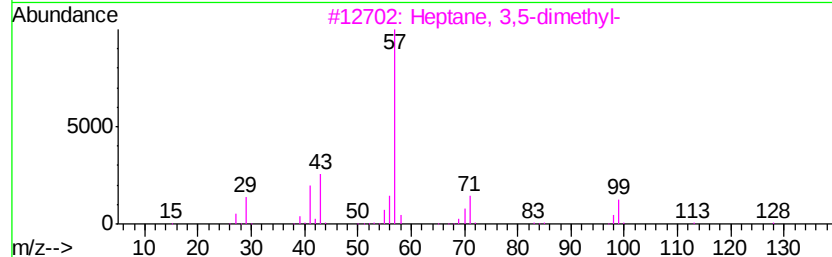
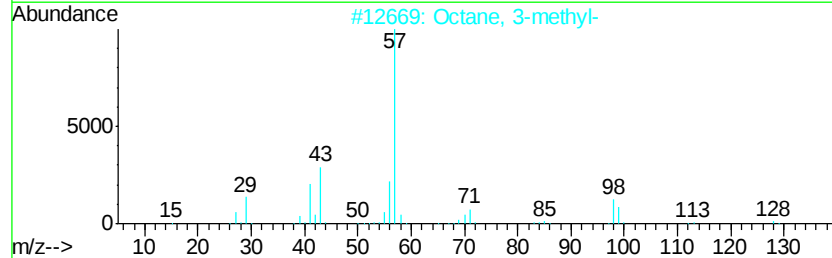
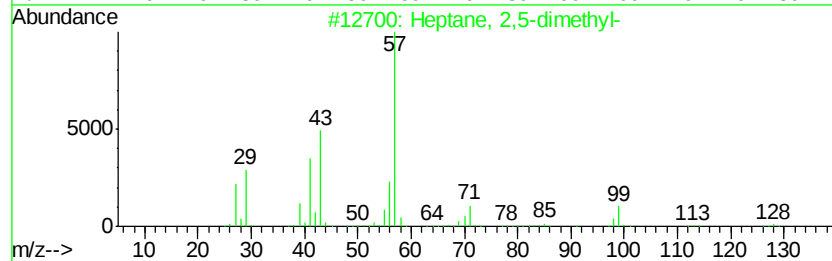
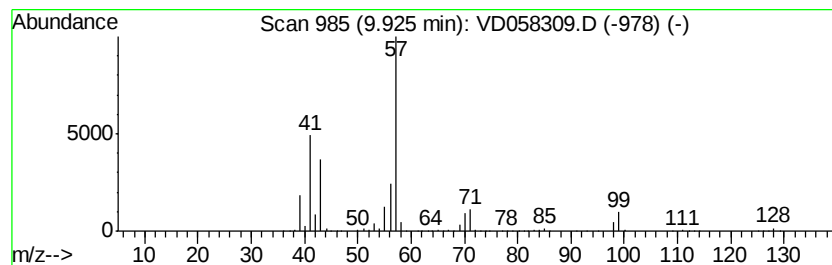
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptane, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	190.68 ug/l	39475400	Chlorobenzene-d5	10.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	91	
2	Octane, 3-methyl-	128	C9H20	002216-33-3	80	
3	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	64	
4	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	53	
5	2,2,4,4-Tetramethyloctane	170	C12H26	062183-79-3	53	



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WP021SB470-27-29-180618

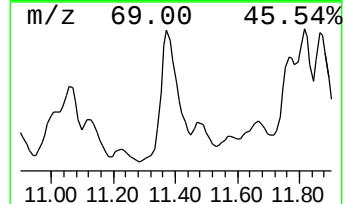
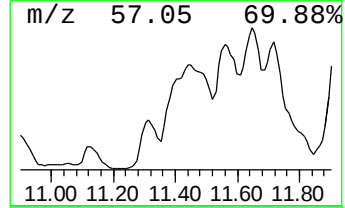
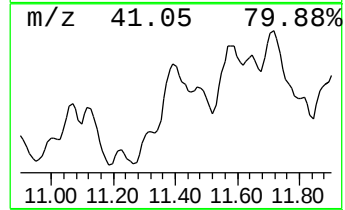
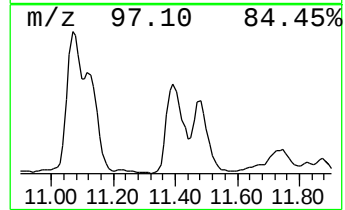
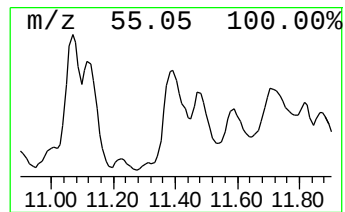
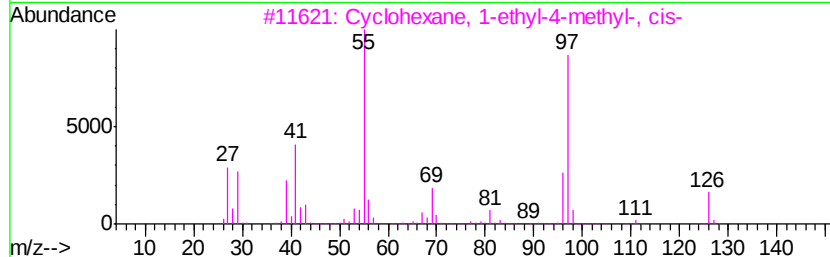
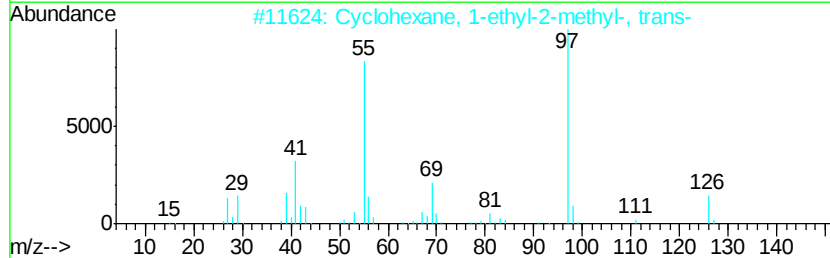
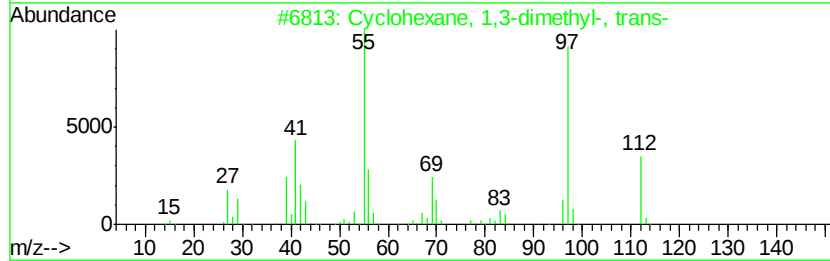
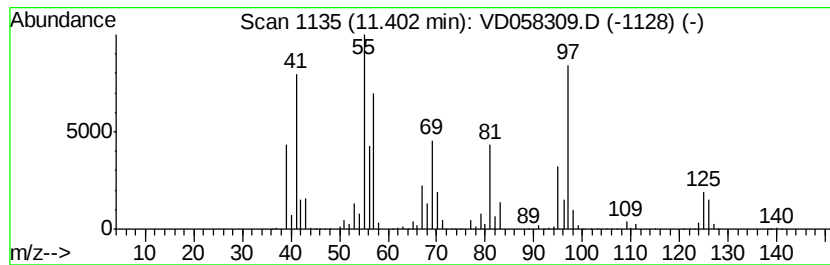
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Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, 1,3-dimethyl-,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	277.79 ug/l	57509400	Chlorobenzene-d5	10.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	58
2		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	46
3		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	46
4		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	46
5		3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062218\
Data File : VD058309.D
Acq On : 22 Jun 2018 21:12
Operator : VA/AP
Sample : J3663-12
Misc : 8.94g/5ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

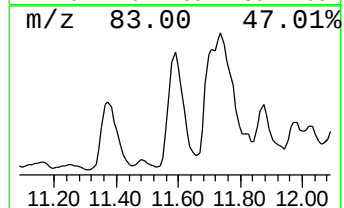
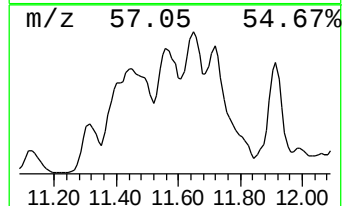
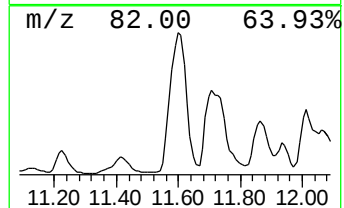
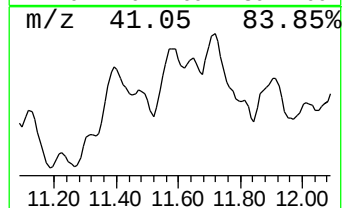
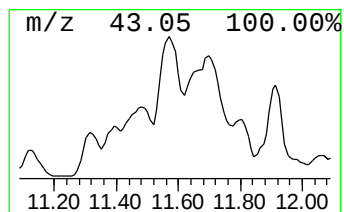
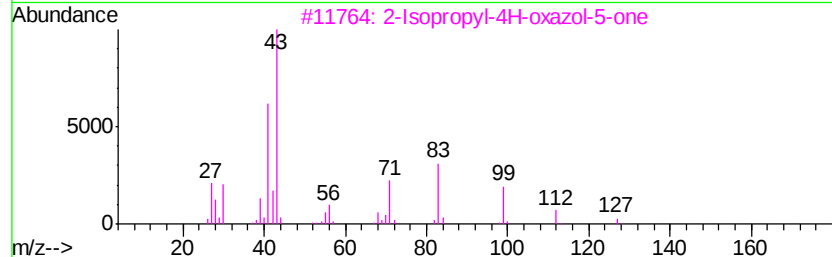
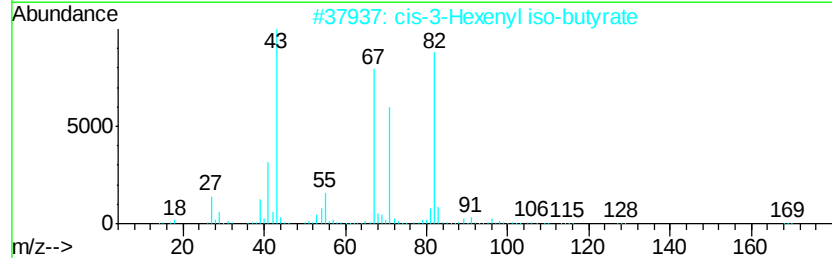
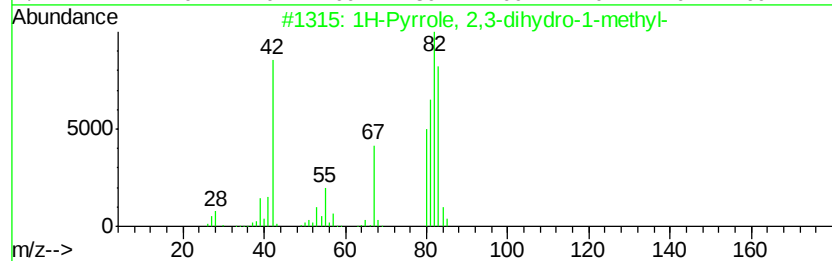
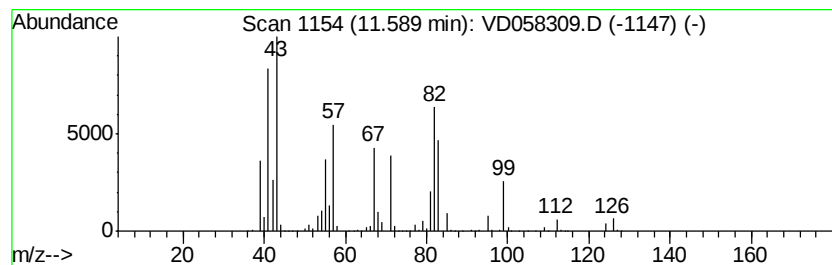
Instrument :
MSVOA_D
ClientSampleID :
WP021SB470-27-29-180618

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 6 unknown11.59 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.59	247.33 ug/l	51202600	Chlorobenzene-d5	10.69	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Pyrrole, 2,3-dihydro-1-methyl-	83	C5H9N	033838-11-8	47
2	cis-3-Hexenyl iso-butyrate	170	C10H18O2	041519-23-7	43
3	2-Isopropyl-4H-oxazol-5-one	127	C6H9NO2	1000190-01-9	38
4	3-Hexen-1-ol, (E)-	100	C6H12O	000928-97-2	22
5	Azeleoneitrile	150	C9H14N2	001675-69-0	22



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062218\
Data File : VD058309.D
Acq On : 22 Jun 2018 21:12
Operator : VA/AP
Sample : J3663-12
Misc : 8.94g/5ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

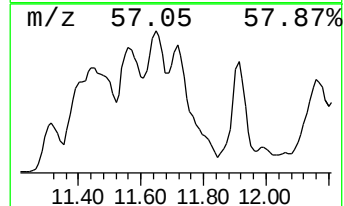
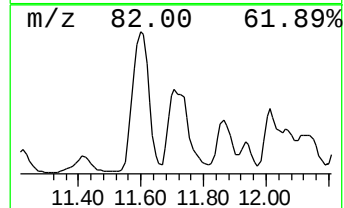
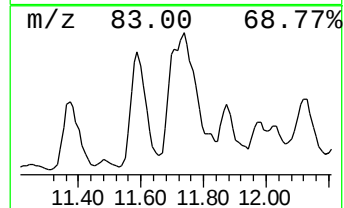
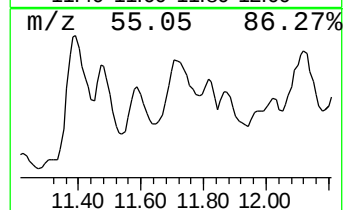
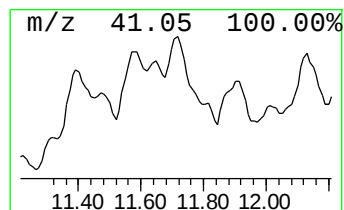
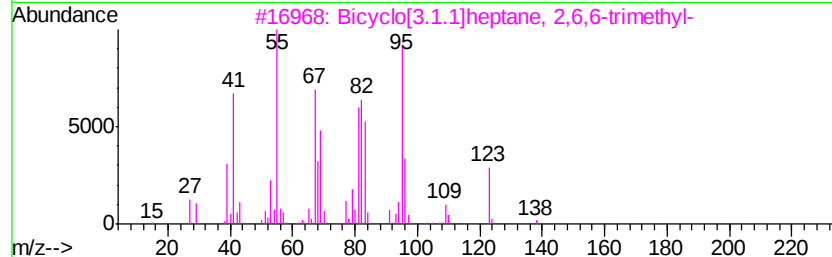
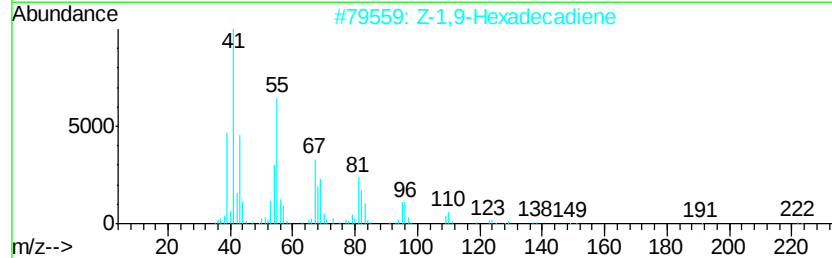
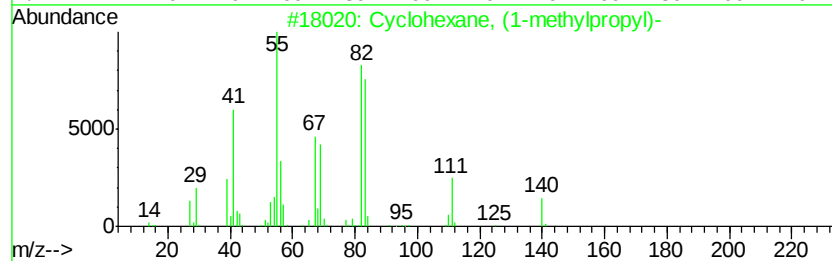
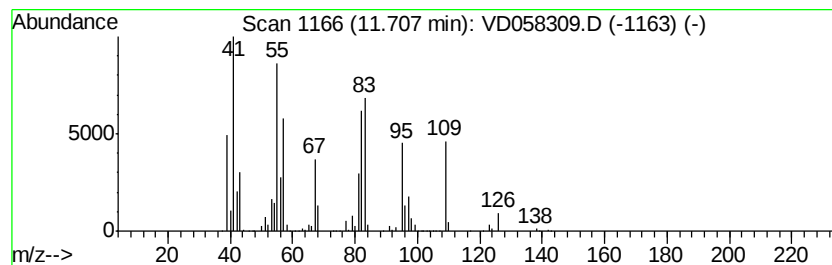
Instrument :
MSVOA_D
ClientSampleId :
WP021SB470-27-29-180618

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 7 unknown11.71 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.71	293.99 ug/l	60864200	Chlorobenzene-d5	10.69		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	46
2		Z-1,9-Hexadecadiene	222	C16H30	1000245-71-3	43
3		Bicyclo[3.1.1]heptane, 2,6,6-tri...	138	C10H18	000473-55-2	38
4		4-t-Butyl-2-(1-methyl-2-nitroeth...	241	C13H23NO3	1000192-74-8	38
5		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062218\
Data File : VD058309.D
Acq On : 22 Jun 2018 21:12
Operator : VA/AP
Sample : J3663-12
Misc : 8.94g/5ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

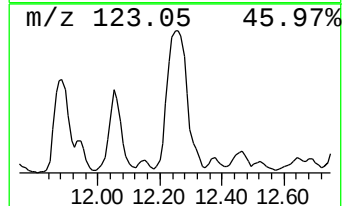
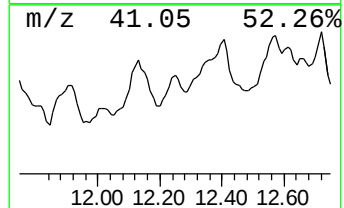
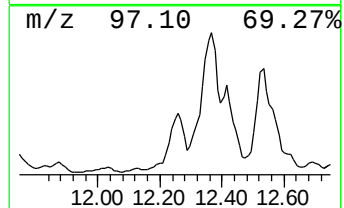
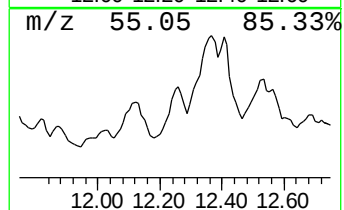
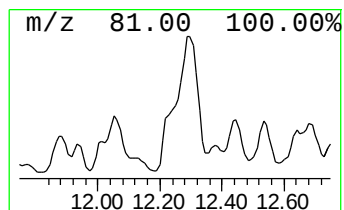
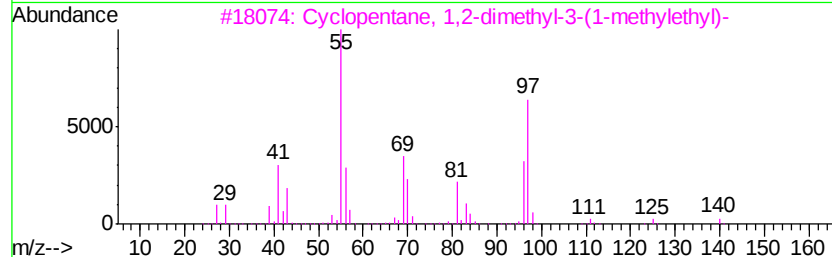
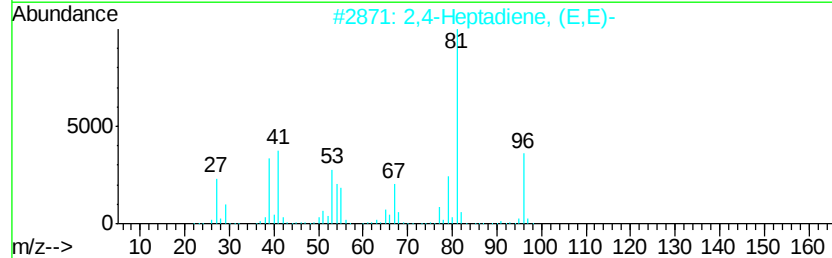
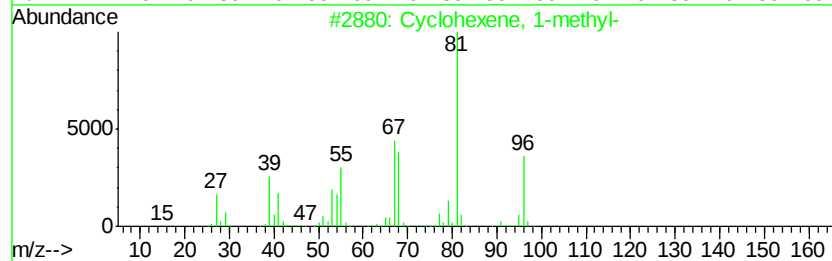
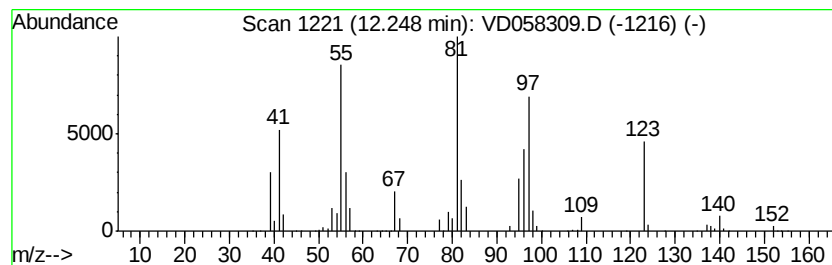
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ClientSampleId :
WP021SB470-27-29-180618

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown12.25 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	163.22 ug/l	19645800	1,4-Dichlorobenzene-d4	12.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexene, 1-methyl-	96 C7H12	000591-49-1 38
2		2,4-Heptadiene, (E,E)-	96 C7H12	002384-94-3 35
3		Cyclopentane, 1,2-dimethyl-3-(1-...	140 C10H20	000489-20-3 35
4		Cyclopentane, 1-methyl-3-(2-meth...	140 C10H20	029053-04-1 35
5		1-Nonylcycloheptane	224 C16H32	1000371-47-7 35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062218\
Data File : VD058309.D
Acq On : 22 Jun 2018 21:12
Operator : VA/AP
Sample : J3663-12
Misc : 8.94g/5ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

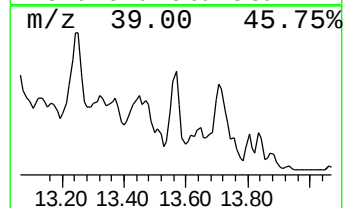
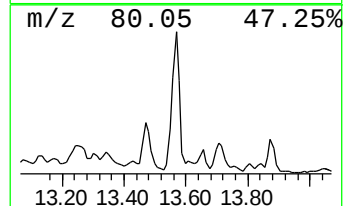
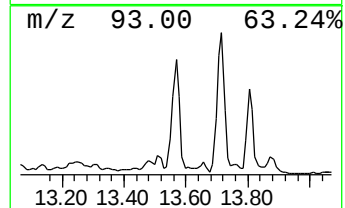
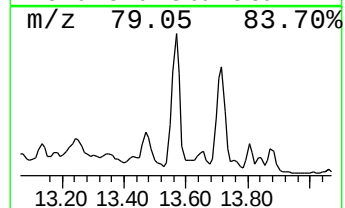
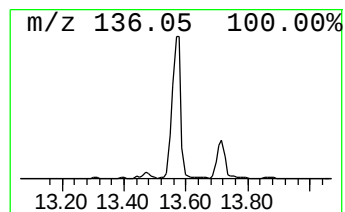
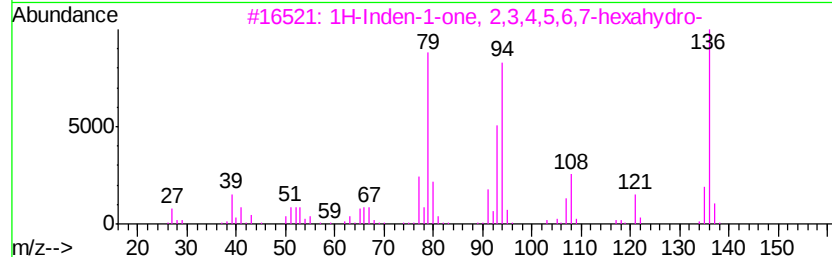
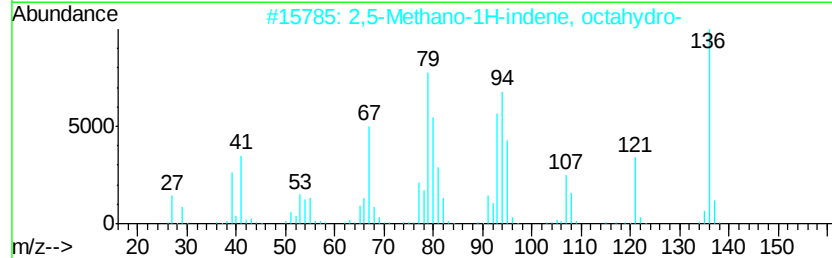
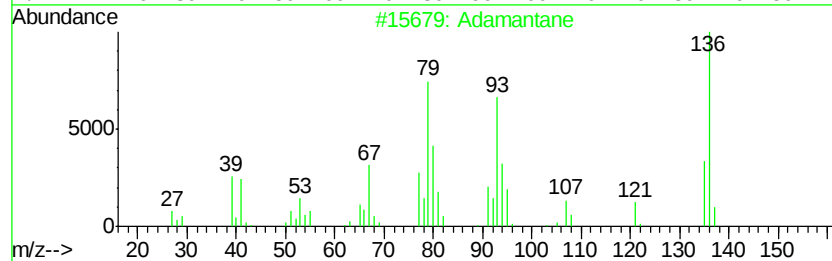
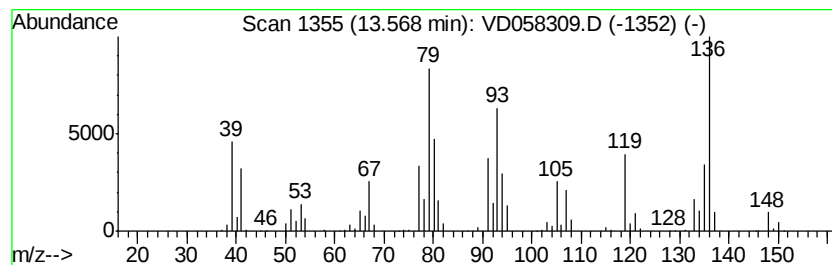
Instrument :
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ClientSampleId :
WP021SB470-27-29-180618

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 9 Adamantane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.57	213.87 ug/l	25742300	1,4-Dichlorobenzene-d4	12.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Adamantane	136 C10H16	000281-23-2	91
2	2,5-Methano-1H-indene, octahydro-	136 C10H16	019026-94-9	64
3	1H-Inden-1-one, 2,3,4,5,6,7-hexa...	136 C9H12O	022118-00-9	58
4	Bicyclo[3.3.1]non-2-en-9-one	136 C9H12O	004844-11-5	58
5	Bicyclo[2.2.1]heptane, 2,2-dimet...	136 C10H16	000497-32-5	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062218\
Data File : VD058309.D
Acq On : 22 Jun 2018 21:12
Operator : VA/AP
Sample : J3663-12
Misc : 8.94g/5ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WP021SB470-27-29-180618

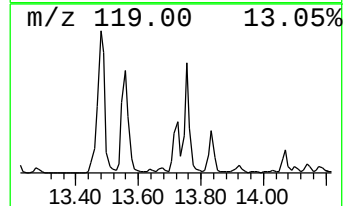
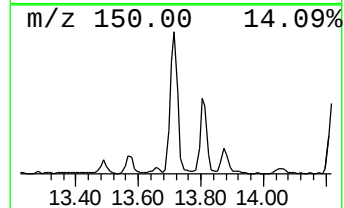
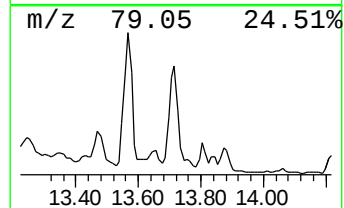
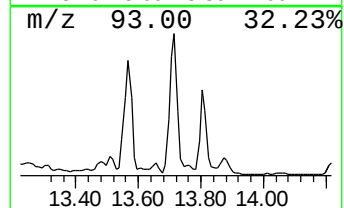
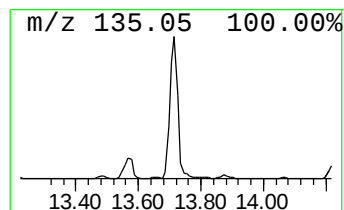
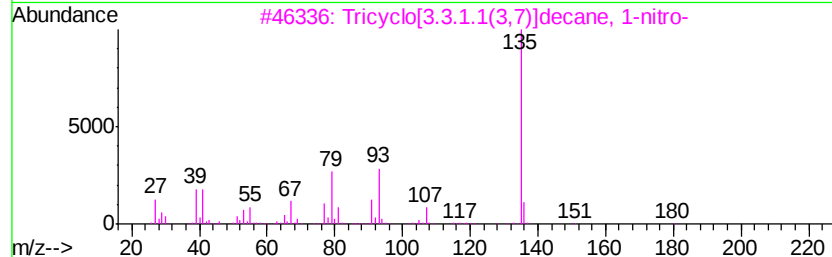
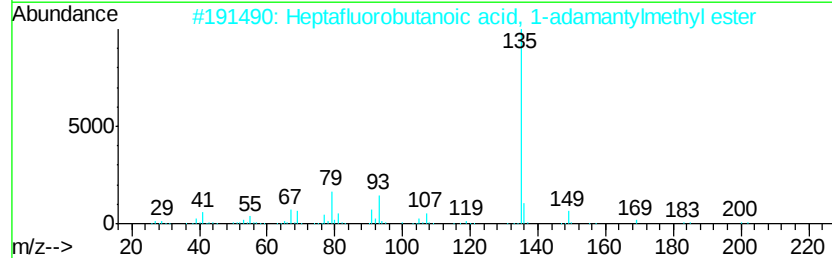
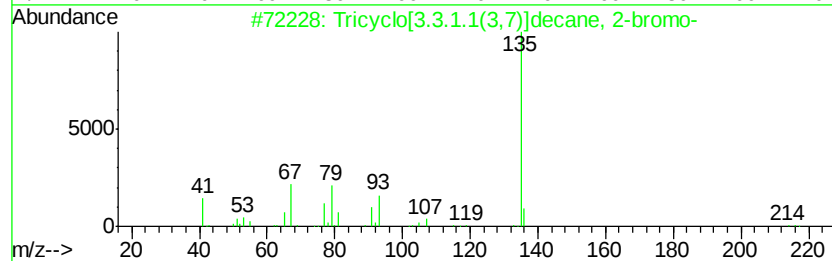
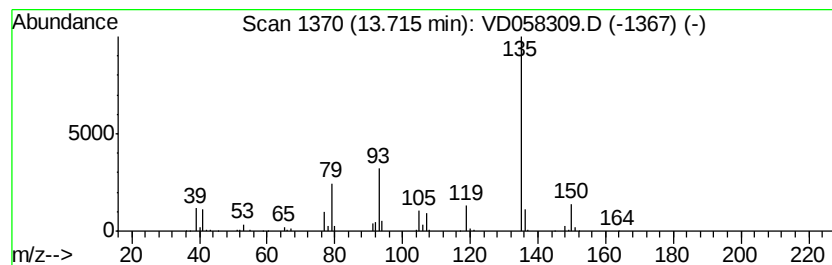
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 10 Tricyclo[3.3.1.1(3,7)]decan... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	346.10 ug/l	41658500	1,4-Dichlorobenzene-d4	12.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[3.3.1.1(3,7)]decane, 2-...	214	C10H15Br	007314-85-4	64
2		Heptafluorobutanoic acid, 1-adam...	362	C15H17F7O2	1000282-96-9	64
3		Tricyclo[3.3.1.1(3,7)]decane, 1-...	181	C10H15NO2	007575-82-8	64
4		1-Adamantanecarboxylic acid chlo...	198	C11H15ClO	002094-72-6	64
5		1-Adamantan-1-yl-6-chloro-7-fluo...	389	C21H21ClFN03	1000210-85-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD062218\
Data File : VD058309.D
Acq On : 22 Jun 2018 21:12
Operator : VA/AP
Sample : J3663-12
Misc : 8.94g/5ml/MSVOA_D/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WP021SB470-27-29-180618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.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
Pentane, 2,2,4-tr...	6.21	406.8	ug/l	18496100	2	6.69	2273620	50.0
Pentane, 2,3,4-tr...	7.91	285.9	ug/l	13001100	2	6.69	2273620	50.0
Pentane, 2,3,3-tr...	8.08	391.8	ug/l	17817300	2	6.69	2273620	50.0
Heptane, 2,5-dime...	9.93	190.7	ug/l	39475400	3	10.69	10351300	50.0
Cyclohexane, 1,3-...	11.40	277.8	ug/l	57509400	3	10.69	10351300	50.0
unknown11.59	11.59	247.3	ug/l	51202600	3	10.69	10351300	50.0
unknown11.71	11.71	294.0	ug/l	60864200	3	10.69	10351300	50.0
unknown12.25	12.25	163.2	ug/l	19645800	4	12.92	6018210	50.0
Adamantane	13.57	213.9	ug/l	25742300	4	12.92	6018210	50.0
Tricyclo[3.3.1.1(...	13.72	346.1	ug/l	41658500	4	12.92	6018210	50.0