

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU061818\
 Data File : VU024690.D
 Acq On : 18 Jun 2018 14:35
 Operator : MD/SY
 Sample : J3463-08
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 T11-MW-9-9.0-061218

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_U\METHOD\82U061318W.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	1.088	136	155	176	rBV	816691	961072	20.07%	2.346%
2	2.323	528	539	551	rBV	57552	90430	1.89%	0.221%
3	2.445	566	577	595	rBV	32235	57573	1.20%	0.141%
4	4.889	1319	1337	1354	rBV2	182562	414014	8.65%	1.011%
5	4.989	1354	1368	1388	rVB	254965	571068	11.93%	1.394%
6	5.313	1456	1469	1485	rBV	178049	375064	7.83%	0.916%
7	5.397	1485	1495	1511	rVB4	22476	50057	1.05%	0.122%
8	5.892	1634	1649	1671	rBV	355691	707526	14.78%	1.727%
9	6.419	1795	1813	1831	rBV	83514	180283	3.77%	0.440%
10	7.570	2149	2171	2184	rBV	684281	1231648	25.72%	3.006%
11	7.641	2184	2193	2207	rVB	88332	156856	3.28%	0.383%
12	7.921	2269	2280	2297	rVB3	25673	48682	1.02%	0.119%
13	8.551	2466	2476	2486	rVB	41324	69034	1.44%	0.169%
14	9.094	2632	2645	2666	rBV	508418	856219	17.88%	2.090%
15	9.377	2716	2733	2749	rBV3	51495	116506	2.43%	0.284%
16	9.783	2853	2859	2883	rVB2	28852	57047	1.19%	0.139%
17	9.956	2890	2913	2925	rBV3	31249	69375	1.45%	0.169%
18	10.049	2930	2942	2960	rBV5	25094	52223	1.09%	0.127%
19	10.168	2969	2979	2990	rBV	108932	170237	3.56%	0.416%
20	10.310	3009	3023	3037	rBV	495753	794707	16.60%	1.940%
21	10.590	3099	3110	3121	rBV	184891	289172	6.04%	0.706%
22	10.709	3133	3147	3157	rBV	43454	78298	1.64%	0.191%
23	10.773	3157	3167	3176	rBV	30693	48026	1.00%	0.117%
24	10.979	3220	3231	3250	rVB4	37889	73789	1.54%	0.180%
25	11.152	3275	3285	3298	rVB	181471	277514	5.80%	0.677%
26	11.294	3312	3329	3333	rBV	60290	103363	2.16%	0.252%
27	11.326	3333	3339	3353	rVB2	153155	240696	5.03%	0.588%
28	11.487	3378	3389	3404	rBV	607529	958834	20.03%	2.340%
29	11.573	3406	3416	3426	rVB	72838	106094	2.22%	0.259%
30	11.692	3442	3453	3460	rBV	42407	64524	1.35%	0.157%
31	11.773	3460	3478	3497	rVV4	370511	1062164	22.18%	2.593%
32	11.869	3497	3508	3530	rVB5	138401	377064	7.88%	0.920%
33	11.985	3531	3544	3557	rBV2	103472	169757	3.55%	0.414%
34	12.149	3586	3595	3600	rBV	43695	64386	1.34%	0.157%

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35	12.181	3600	3605	3612	rVV	38823	52983	1.11%	0.129%
36	12.255	3613	3628	3635	rVV	515571	807904	16.87%	1.972%
37	12.290	3635	3639	3649	rVB2	132356	182918	3.82%	0.446%
38	12.358	3650	3660	3681	rBV4	298808	736743	15.39%	1.798%
39	12.541	3709	3717	3731	rVB2	113026	192869	4.03%	0.471%
40	12.654	3731	3752	3761	rBV	392966	721536	15.07%	1.761%
41	12.705	3761	3768	3784	rVB2	108204	201516	4.21%	0.492%
42	12.789	3784	3794	3809	rBV3	117422	185236	3.87%	0.452%
43	12.882	3810	3823	3829	rBV2	82169	135215	2.82%	0.330%
44	12.930	3830	3838	3843	rVV4	89833	146181	3.05%	0.357%
45	12.966	3844	3849	3857	rVB	170616	230994	4.82%	0.564%
46	13.123	3878	3898	3908	rBV2	1163028	1953394	40.80%	4.768%
47	13.242	3928	3935	3941	rBV	72609	106993	2.23%	0.261%
48	13.294	3942	3951	3971	rVB4	199500	382759	7.99%	0.934%
49	13.409	3978	3987	3994	rBV2	89139	150192	3.14%	0.367%
50	13.461	3995	4003	4010	rVB	133737	202910	4.24%	0.495%
51	13.512	4010	4019	4027	rBV3	318949	473371	9.89%	1.155%
52	13.580	4027	4040	4044	rBV2	116010	216634	4.52%	0.529%
53	13.612	4045	4050	4057	rVB	125819	161939	3.38%	0.395%
54	13.744	4076	4091	4102	rBV	2949325	4787920	100.00%	11.687%
55	13.789	4102	4105	4116	rVV	127808	177400	3.71%	0.433%
56	13.860	4116	4127	4140	rVV2	244292	419338	8.76%	1.024%
57	13.956	4140	4157	4169	rVV3	225441	541228	11.30%	1.321%
58	14.017	4169	4176	4185	rVV2	107561	186952	3.90%	0.456%
59	14.065	4186	4191	4207	rVB5	44404	74585	1.56%	0.182%
60	14.155	4207	4219	4236	rBV2	146487	320269	6.69%	0.782%
61	14.284	4245	4259	4265	rBV8	25736	55803	1.17%	0.136%
62	14.351	4265	4280	4300	rVV4	486326	1072861	22.41%	2.619%
63	14.480	4312	4320	4330	rVV2	116952	207558	4.34%	0.507%
64	14.544	4331	4340	4352	rVB4	165454	304823	6.37%	0.744%
65	14.609	4352	4360	4370	rVB4	58176	92078	1.92%	0.225%
66	14.679	4371	4382	4399	rBV2	345962	640356	13.37%	1.563%
67	14.821	4418	4426	4432	rBV2	46857	72656	1.52%	0.177%
68	14.879	4432	4444	4455	rVV	2433697	3877142	80.98%	9.464%
69	14.933	4455	4461	4474	rVB5	112849	213007	4.45%	0.520%
70	15.011	4477	4485	4490	rBV4	38235	54844	1.15%	0.134%
71	15.065	4490	4502	4516	rBV	2764701	4330930	90.46%	10.572%

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72	15.130	4516	4522	4529	rBV3	40345	59623	1.25%	0.146%
73	15.326	4578	4583	4599	rVB10	18669	47945	1.00%	0.117%
74	15.593	4653	4666	4682	rVB5	106490	289066	6.04%	0.706%
75	15.676	4683	4692	4702	rVV4	31384	49070	1.02%	0.120%
76	15.802	4716	4731	4738	rBV	200199	348658	7.28%	0.851%
77	15.917	4753	4767	4781	rBV	537265	963607	20.13%	2.352%
78	16.062	4801	4812	4819	rBV	825384	1324212	27.66%	3.232%
79	16.101	4819	4824	4839	rVB	475073	692271	14.46%	1.690%
80	16.181	4839	4849	4862	rBV3	42601	80928	1.69%	0.198%
81	16.265	4864	4875	4876	rBV2	136283	155630	3.25%	0.380%
82	16.290	4878	4883	4904	rVB2	224836	392716	8.20%	0.959%
83	16.435	4916	4928	4941	rBV	176065	277652	5.80%	0.678%
84	16.740	5012	5023	5040	rBV	328199	545142	11.39%	1.331%
85	17.030	5104	5113	5128	rVB2	30581	74496	1.56%	0.182%
86	17.371	5203	5219	5234	rVB4	17290	51279	1.07%	0.125%

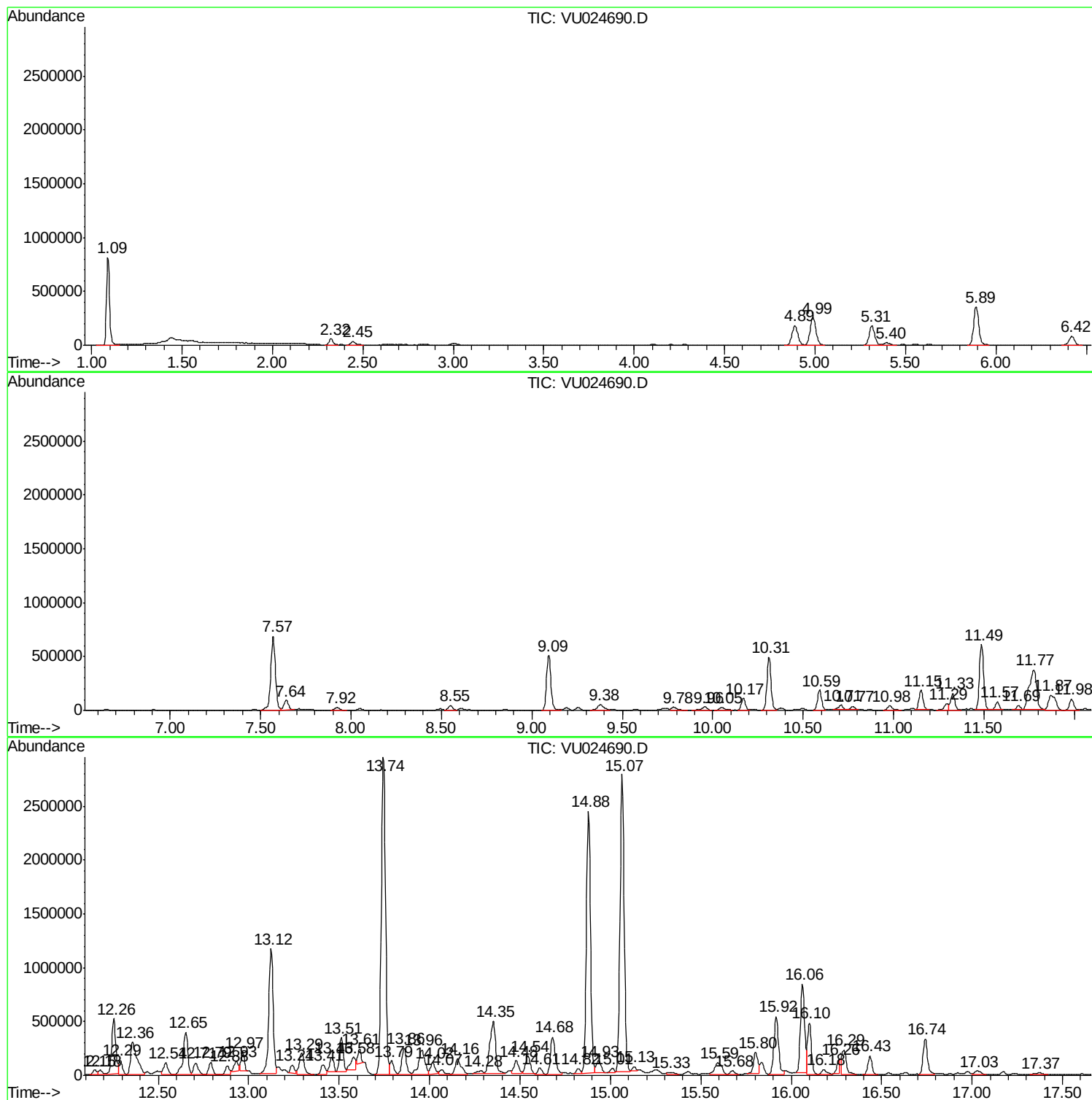
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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



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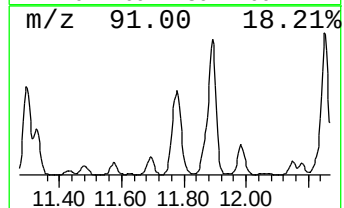
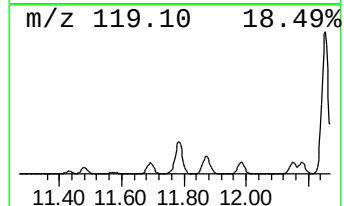
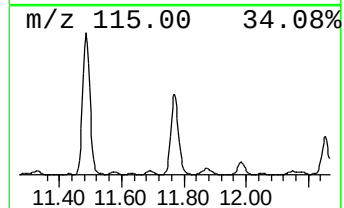
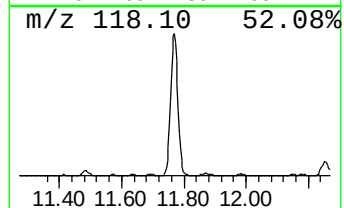
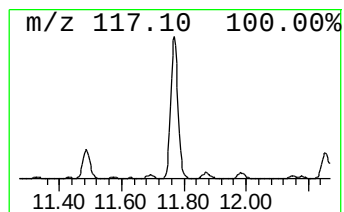
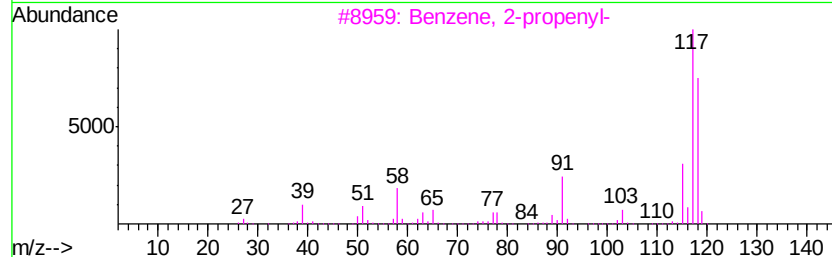
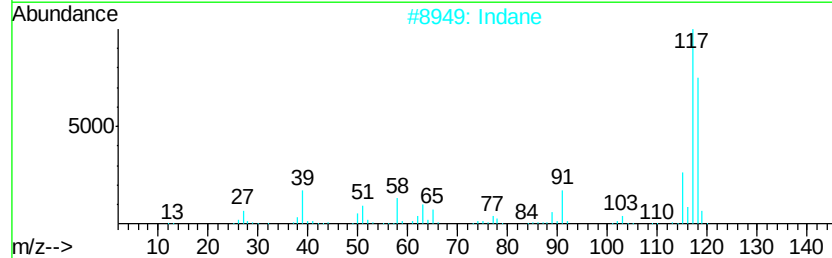
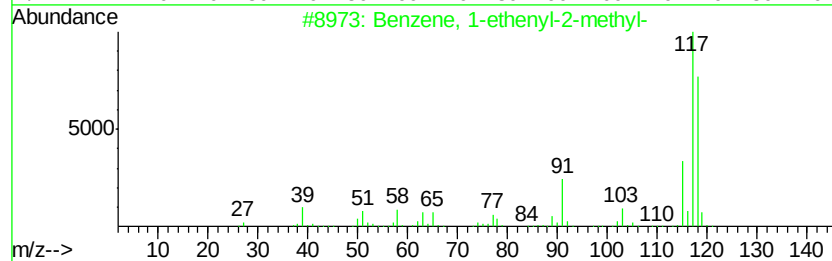
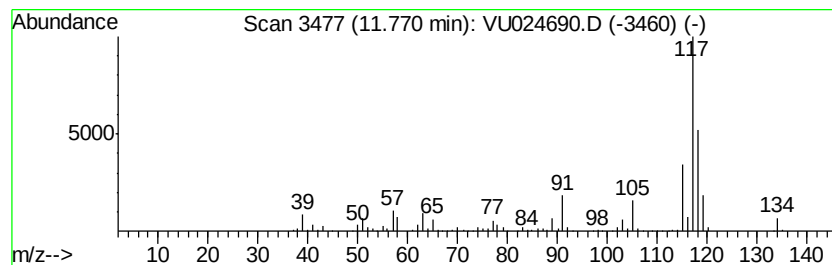
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethenyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.77	55.39 ug/l	1062160	1,4-Dichlorobenzene-d4	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
2	Indane	118	C9H10	000496-11-7	70
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	64



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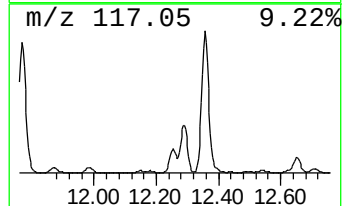
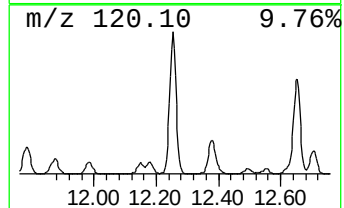
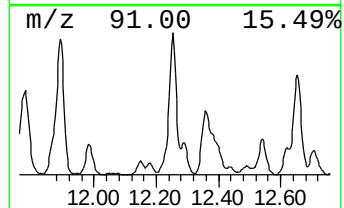
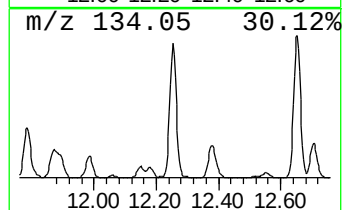
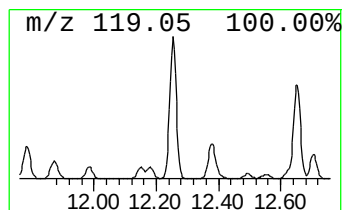
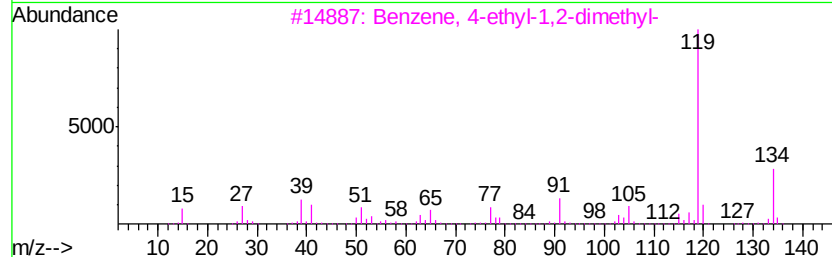
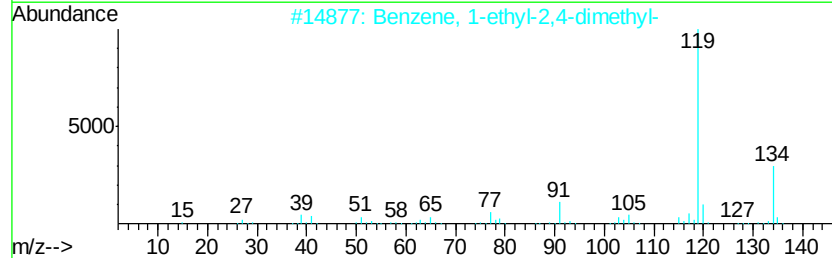
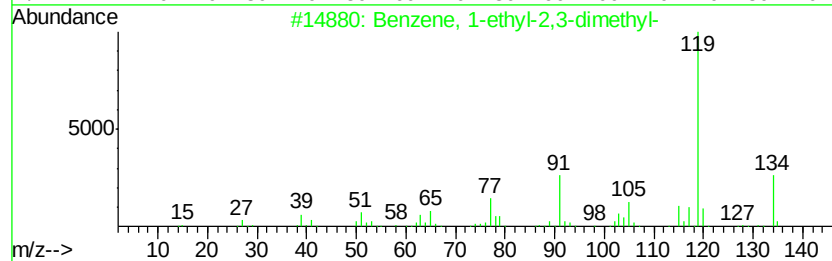
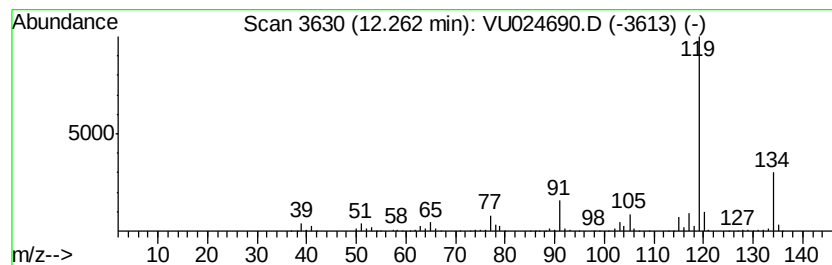
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	42.13 ug/l	807904	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5		o-Cymene	134	C10H14	000527-84-4	94



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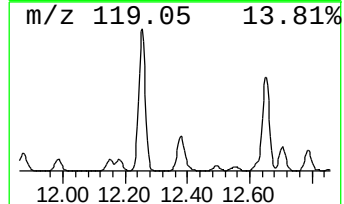
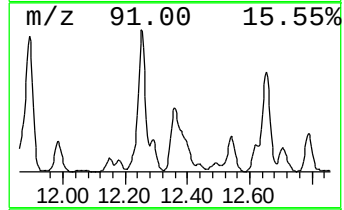
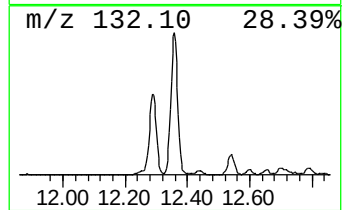
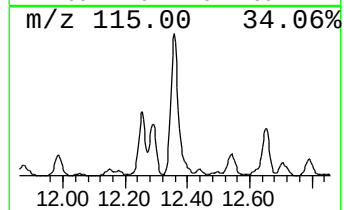
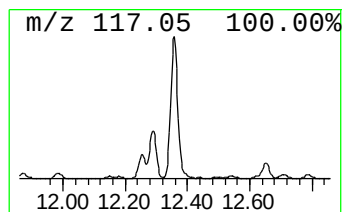
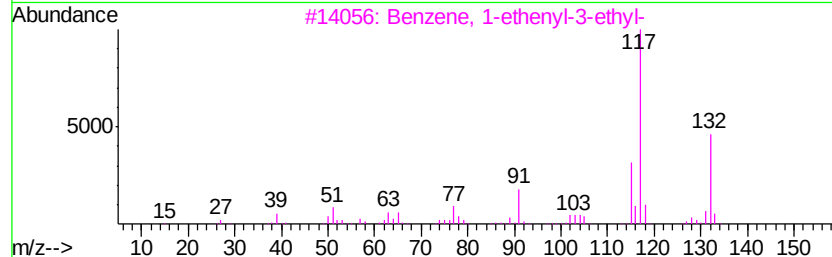
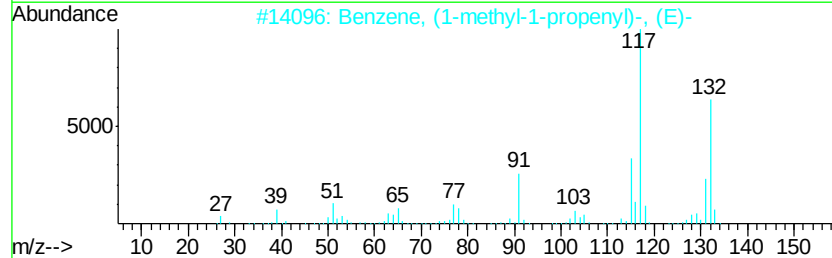
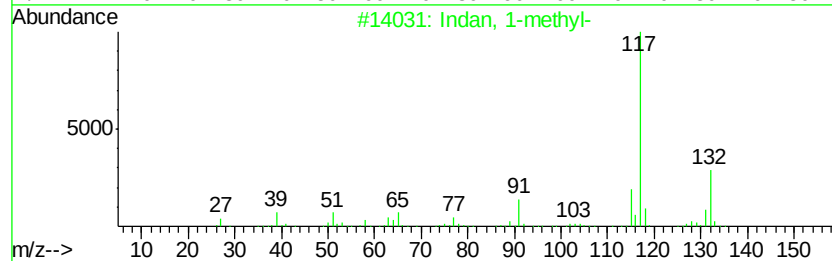
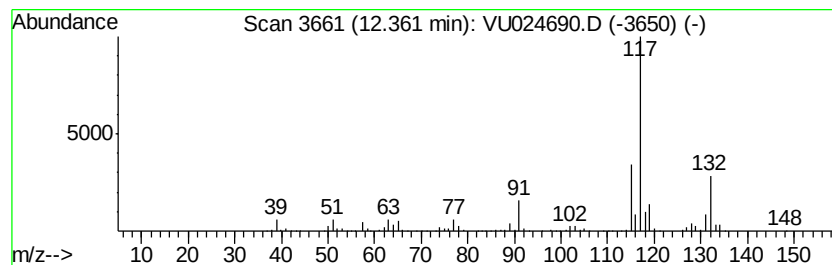
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Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	38.42 ug/l	736743	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	87
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
5		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83



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T11-MW-9-9.0-061218

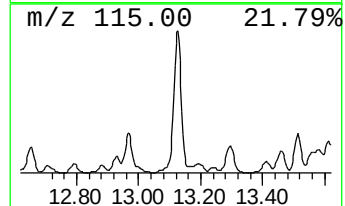
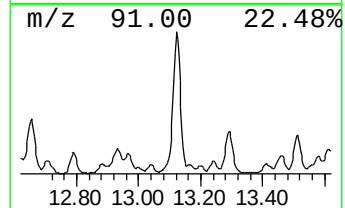
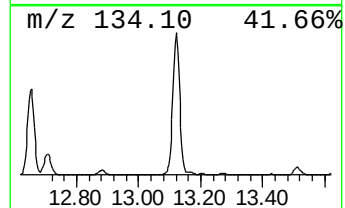
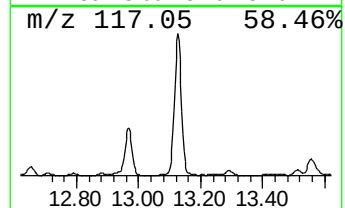
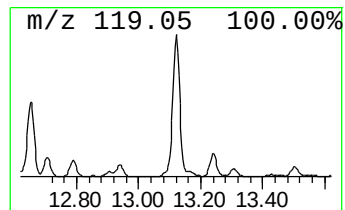
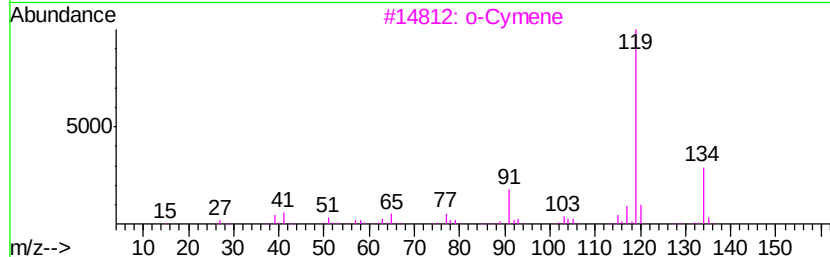
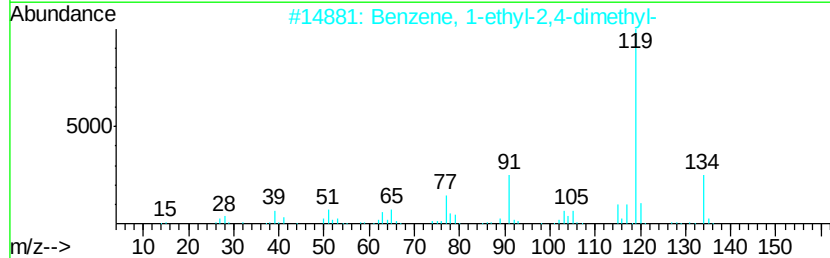
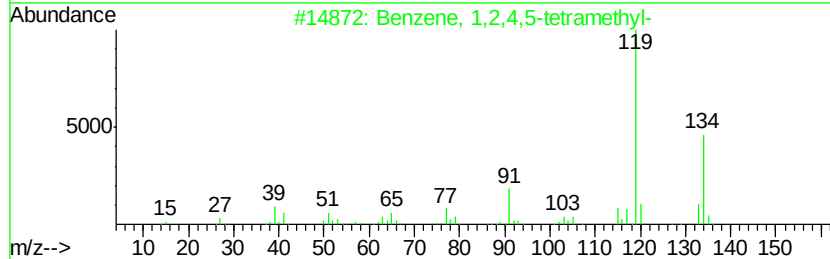
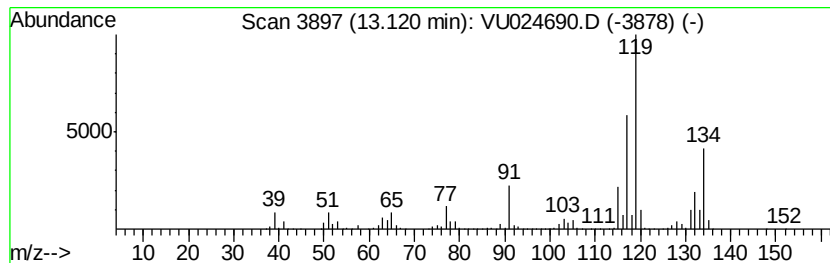
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	101.86 ug/l	1953390	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
2		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	60
3		o-Cymene	134	C10H14	000527-84-4	60
4		p-Cymene	134	C10H14	000099-87-6	60
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024690.D
Acq On : 18 Jun 2018 14:35
Operator : MD/SY
Sample : J3463-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 12 Sample Multiplier: 1

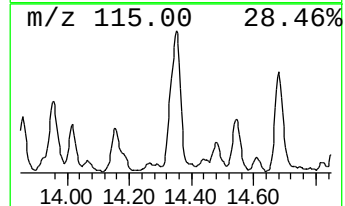
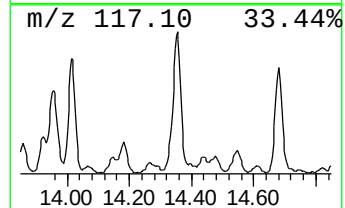
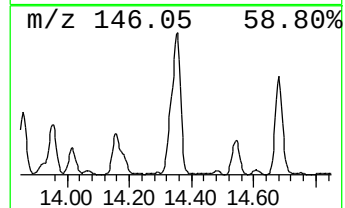
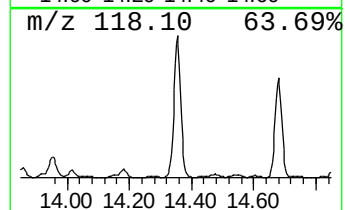
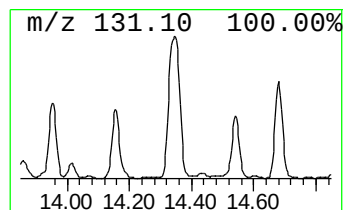
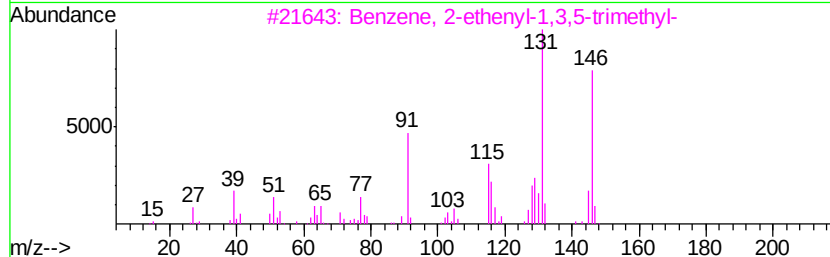
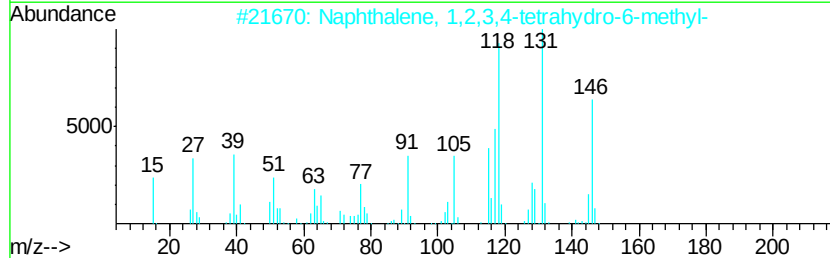
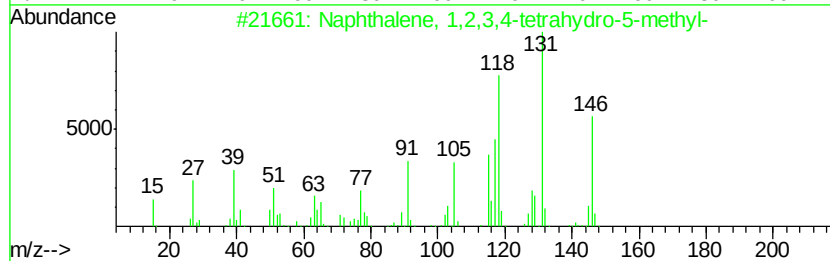
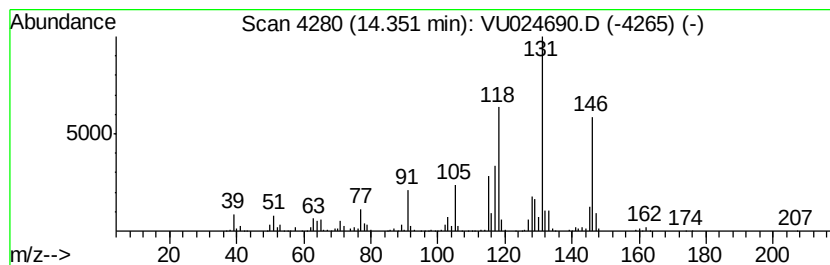
Instrument :
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ClientSampleId :
T11-MW-9-9.0-061218

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.35	55.95 ug/l	1072860	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 97
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 96
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 86
4		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 70
5		Benzene, (1,1-dimethyl-2-propenyl)-	146 C11H14	018321-36-3 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024690.D
Acq On : 18 Jun 2018 14:35
Operator : MD/SY
Sample : J3463-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 12 Sample Multiplier: 1

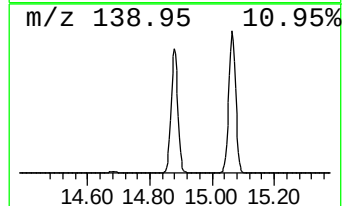
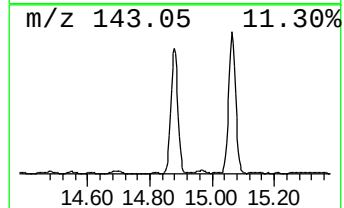
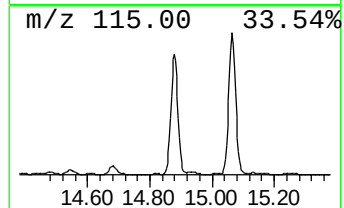
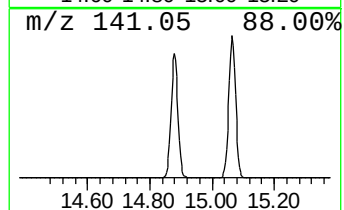
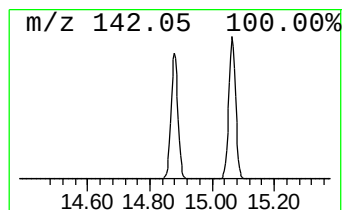
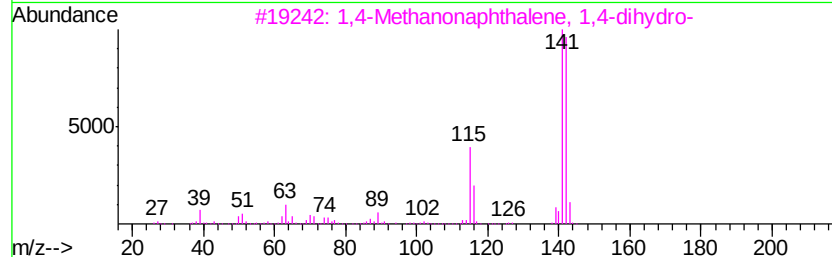
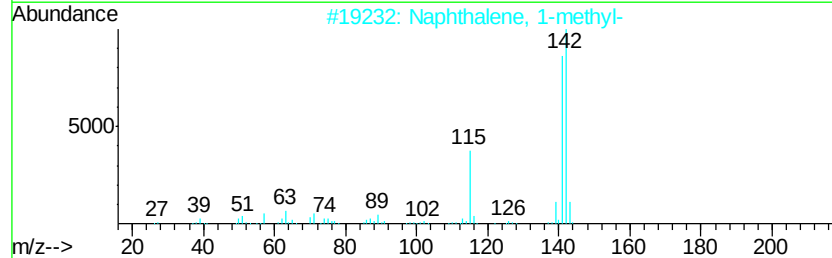
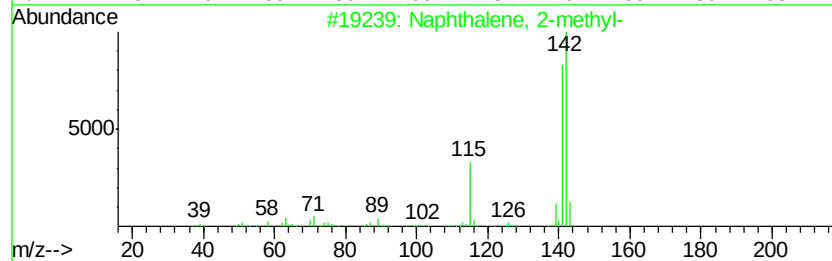
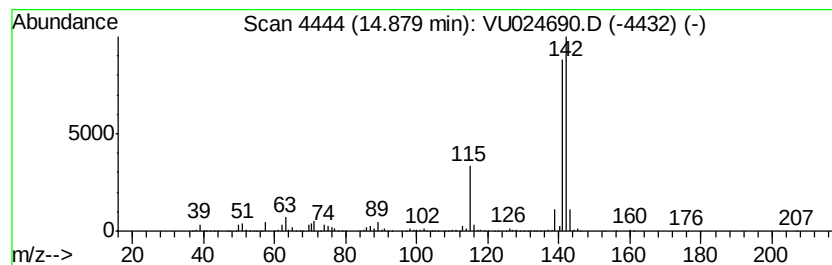
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ClientSampleId :
T11-MW-9-9.0-061218

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.88	202.18 ug/l	3877140	1,4-Dichlorobenzene-d4	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	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:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024690.D
Acq On : 18 Jun 2018 14:35
Operator : MD/SY
Sample : J3463-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 12 Sample Multiplier: 1

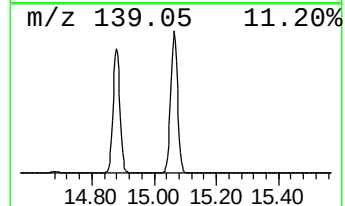
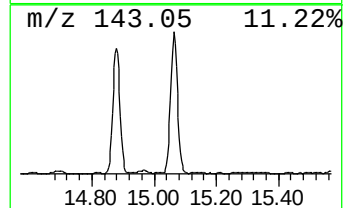
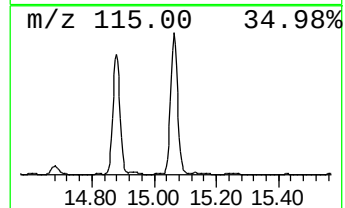
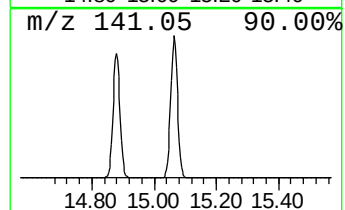
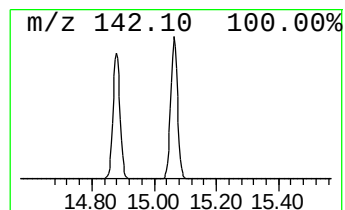
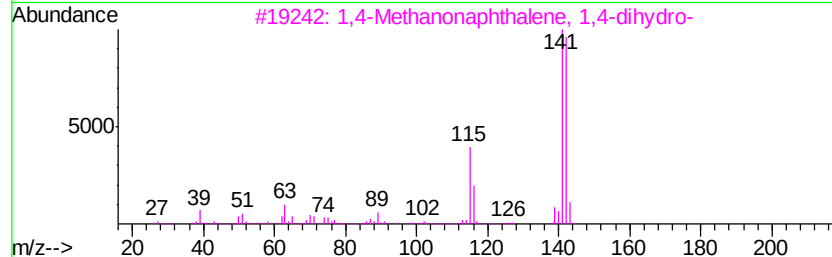
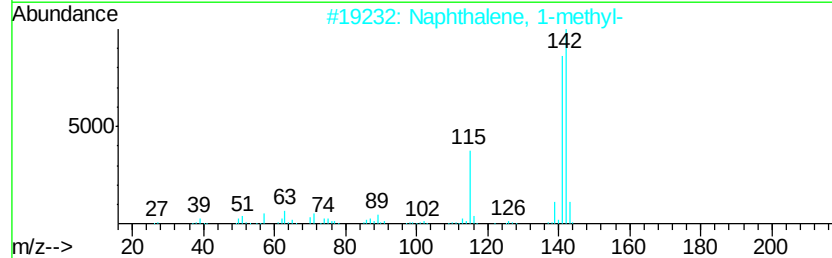
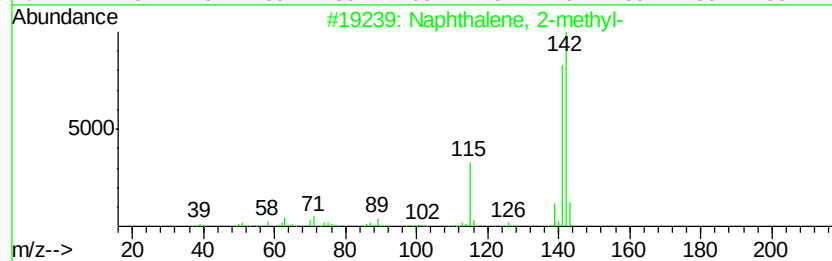
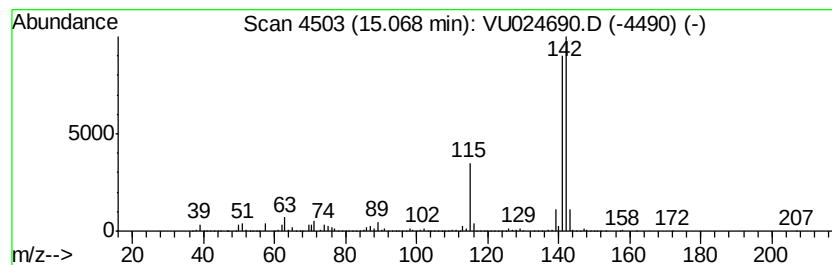
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 8 1,4-Methanonaphthalene, 1,4... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.07	225.84 ug/l	4330930	1,4-Dichlorobenzene-d4	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	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:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024690.D
Acq On : 18 Jun 2018 14:35
Operator : MD/SY
Sample : J3463-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 12 Sample Multiplier: 1

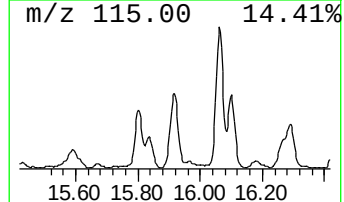
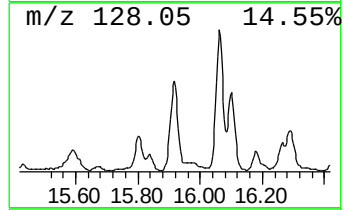
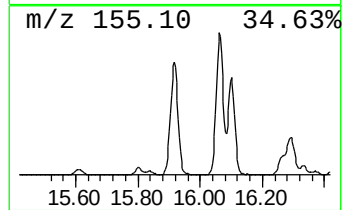
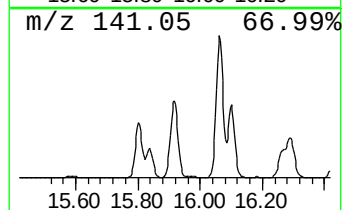
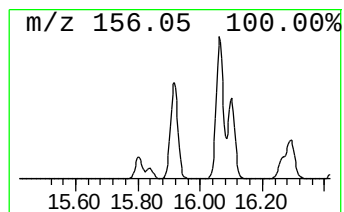
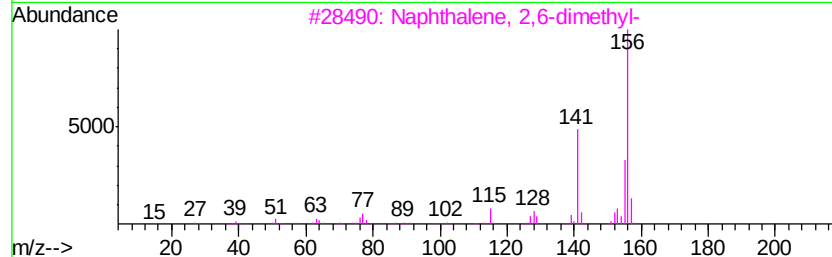
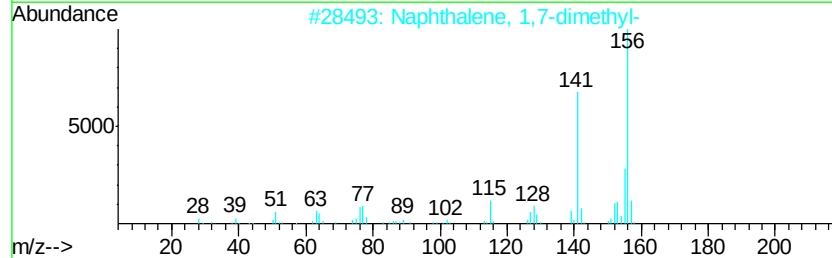
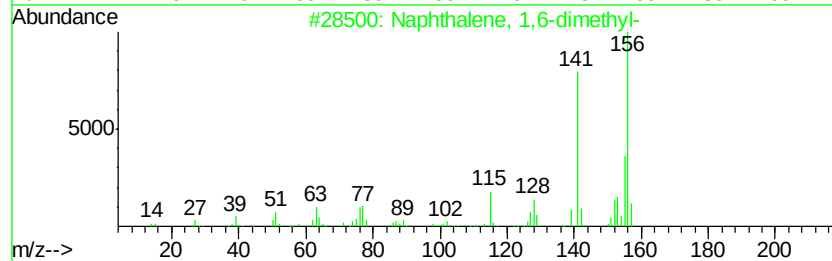
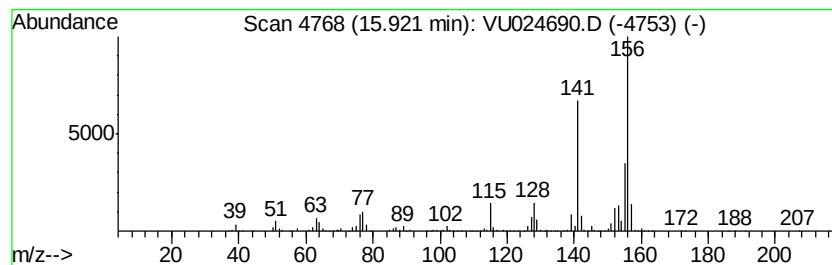
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	50.25 ug/l	963607	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
2		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
3		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
4		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 97
5		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024690.D
Acq On : 18 Jun 2018 14:35
Operator : MD/SY
Sample : J3463-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 12 Sample Multiplier: 1

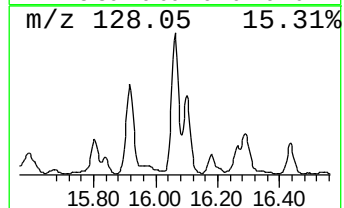
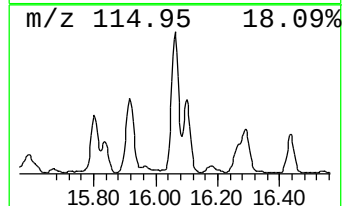
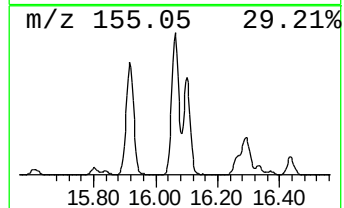
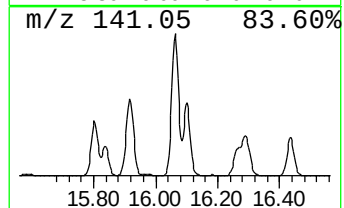
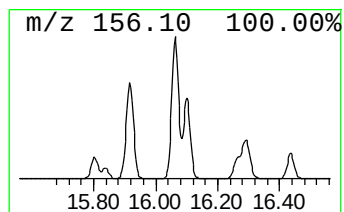
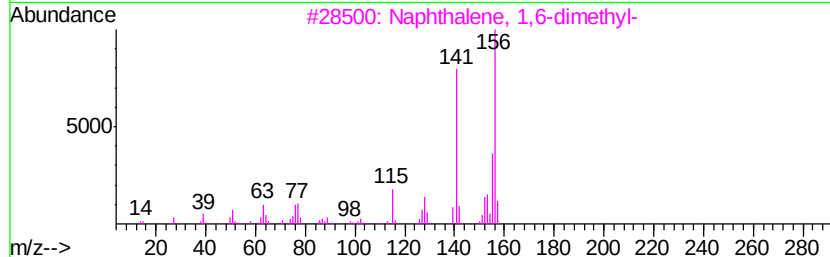
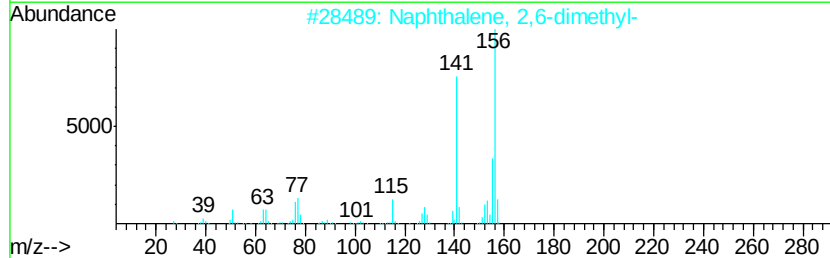
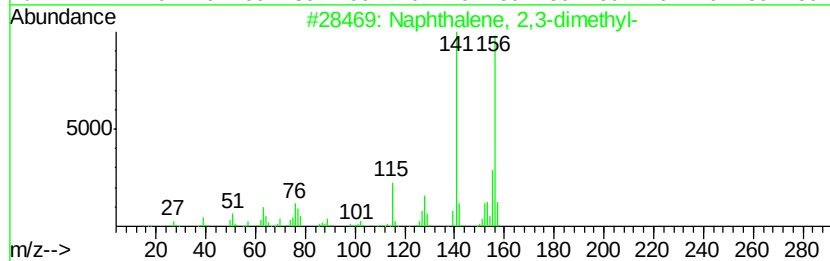
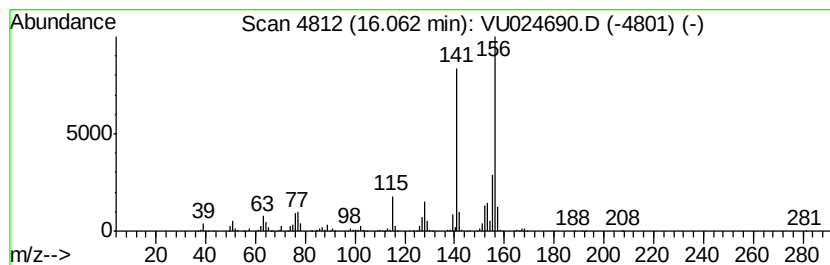
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.06	69.05 ug/l	1324210	1,4-Dichlorobenzene-d4	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061818\
Data File : VU024690.D
Acq On : 18 Jun 2018 14:35
Operator : MD/SY
Sample : J3463-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
T11-MW-9-9.0-061218

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.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
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Benzene, 1-ethyl-...	12.26	42.1	ug/l	807904	4	11.49	958834	50.0
Indan, 1-methyl-	12.36	38.4	ug/l	736743	4	11.49	958834	50.0
Benzene, 1,2,4,5-...	13.12	101.9	ug/l	1953390	4	11.49	958834	50.0
Naphthalene, 1,2,...	14.35	56.0	ug/l	1072860	4	11.49	958834	50.0
Naphthalene, 1-me...	14.88	202.2	ug/l	3877140	4	11.49	958834	50.0
1,4-Methanonaphth...	15.07	225.8	ug/l	4330930	4	11.49	958834	50.0
Naphthalene, 1,6-...	15.92	50.3	ug/l	963607	4	11.49	958834	50.0
Naphthalene, 2,3-...	16.06	69.1	ug/l	1324210	4	11.49	958834	50.0