

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102318\
 Data File : VX005522.D
 Acq On : 23 Oct 2018 15:47
 Operator : JC/MD
 Sample : J5622-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KG-V23608

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.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	2.446	205	212	220	rBV	26005	41596	4.19%	0.440%
2	2.861	274	280	286	rVV	8457	15665	1.58%	0.166%
3	4.696	575	581	588	rBV2	5911	17000	1.71%	0.180%
4	5.507	700	714	725	rBV3	135081	384236	38.73%	4.062%
5	5.665	727	740	755	rBV2	221429	620145	62.51%	6.556%
6	6.068	795	806	829	rVB	138732	379187	38.22%	4.008%
7	6.464	860	871	887	rBV	83325	214361	21.61%	2.266%
8	6.860	925	936	950	rBV	360050	816594	82.32%	8.632%
9	7.884	1097	1104	1114	rBV	69619	123066	12.41%	1.301%
10	8.622	1219	1225	1234	rBV	5175	10363	1.04%	0.110%
11	8.720	1234	1241	1264	rVV	588153	991995	100.00%	10.486%
12	9.408	1350	1354	1363	rVB2	11906	19845	2.00%	0.210%
13	9.549	1372	1377	1382	rBV	46186	62920	6.34%	0.665%
14	9.975	1440	1447	1458	rBV	60246	82976	8.36%	0.877%
15	10.116	1463	1470	1487	rVV	607444	854027	86.09%	9.028%
16	11.140	1632	1638	1651	rBV	496876	642145	64.73%	6.788%
17	12.079	1786	1792	1800	rBV	651381	800102	80.66%	8.458%
18	12.304	1821	1829	1836	rBV2	16244	24264	2.45%	0.256%
19	12.390	1836	1843	1850	rVB6	4489	10286	1.04%	0.109%
20	12.463	1850	1855	1862	rBV2	7039	11033	1.11%	0.117%
21	12.670	1884	1889	1892	rBV	35800	42948	4.33%	0.454%
22	12.701	1892	1894	1899	rVB	9914	13071	1.32%	0.138%
23	12.755	1899	1903	1911	rBV4	29031	47693	4.81%	0.504%
24	12.896	1923	1926	1931	rVB	8212	11478	1.16%	0.121%
25	12.981	1931	1940	1944	rBV	39504	57781	5.82%	0.611%
26	13.085	1952	1957	1961	rBV2	11283	16994	1.71%	0.180%
27	13.152	1964	1968	1971	rBV	9952	11638	1.17%	0.123%
28	13.194	1971	1975	1978	rBV2	16276	22021	2.22%	0.233%
29	13.225	1978	1980	1984	rVB	19650	20771	2.09%	0.220%
30	13.316	1991	1995	1996	rBV3	9685	12378	1.25%	0.131%
31	13.347	1996	2000	2007	rVV2	95159	145976	14.72%	1.543%
32	13.426	2007	2013	2019	rVV2	16187	28449	2.87%	0.301%
33	13.481	2019	2022	2029	rVB2	9029	13694	1.38%	0.145%
34	13.566	2029	2036	2038	rBV	15331	22136	2.23%	0.234%

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Integrator: RTE
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35	13.603	2038	2042	2045	rVV	35430	49094	4.95%	0.519%
36	13.639	2045	2048	2052	rVV2	49121	71074	7.16%	0.751%
37	13.700	2052	2058	2059	rVV3	27558	46570	4.69%	0.492%
38	13.719	2059	2061	2070	rVB2	56875	90975	9.17%	0.962%
39	13.810	2070	2076	2078	rBV2	12838	18098	1.82%	0.191%
40	13.853	2078	2083	2088	rVB4	14622	27675	2.79%	0.293%
41	13.914	2088	2093	2097	rVB	47737	59825	6.03%	0.632%
42	13.987	2097	2105	2109	rBV2	70841	101027	10.18%	1.068%
43	14.030	2109	2112	2116	rVV	25284	30065	3.03%	0.318%
44	14.133	2124	2129	2133	rBV	52470	71085	7.17%	0.751%
45	14.280	2146	2153	2161	rBV3	80689	181857	18.33%	1.922%
46	14.383	2166	2170	2174	rVV	38649	53409	5.38%	0.565%
47	14.432	2174	2178	2182	rVV	69941	88113	8.88%	0.931%
48	14.475	2182	2185	2190	rVB	15546	21225	2.14%	0.224%
49	14.542	2190	2196	2200	rBV2	111740	148849	15.01%	1.573%
50	14.676	2208	2218	2221	rBV3	54583	90625	9.14%	0.958%
51	14.731	2224	2227	2231	rVB	46253	58751	5.92%	0.621%
52	14.786	2232	2236	2241	rBV3	19041	23869	2.41%	0.252%
53	14.840	2241	2245	2248	rBV2	31704	50368	5.08%	0.532%
54	14.877	2248	2251	2254	rVV2	28208	42850	4.32%	0.453%
55	14.907	2254	2256	2260	rVB3	20904	23927	2.41%	0.253%
56	14.968	2260	2266	2272	rBV3	20199	42690	4.30%	0.451%
57	15.115	2282	2290	2294	rBV2	13075	21023	2.12%	0.222%
58	15.249	2307	2312	2320	rVV	60334	105365	10.62%	1.114%
59	15.328	2320	2325	2331	rVB2	16584	30208	3.05%	0.319%
60	15.444	2338	2344	2347	rBV3	21464	35283	3.56%	0.373%
61	15.481	2347	2350	2354	rVV	35089	45071	4.54%	0.476%
62	15.554	2354	2362	2366	rVV	113626	183248	18.47%	1.937%
63	15.596	2366	2369	2374	rVV5	21394	40508	4.08%	0.428%
64	15.688	2378	2384	2388	rVV	278732	441424	44.50%	4.666%
65	15.724	2388	2390	2398	rVB	67259	101453	10.23%	1.072%
66	15.810	2398	2404	2410	rBV5	11541	26326	2.65%	0.278%
67	15.919	2411	2422	2438	rVB	72850	206719	20.84%	2.185%
68	16.072	2441	2447	2458	rVB	43654	84071	8.47%	0.889%
69	16.176	2458	2464	2468	rBV7	8138	14980	1.51%	0.158%
70	16.285	2479	2482	2487	rVB5	7722	12411	1.25%	0.131%
71	16.352	2488	2493	2496	rBV3	6393	11552	1.16%	0.122%

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ClientSampleId :
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Integration Parameters: RTEINT.P

Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	16.419	2501	2504	2505	rBV	10145	11231	1.13%	0.119%
73	16.572	2525	2529	2534	rBV3	8001	13440	1.35%	0.142%
74	16.657	2534	2543	2545	rBV8	10279	24066	2.43%	0.254%
75	16.694	2545	2549	2554	rVB2	21159	38106	3.84%	0.403%
76	16.755	2554	2559	2564	rBV	18114	32525	3.28%	0.344%

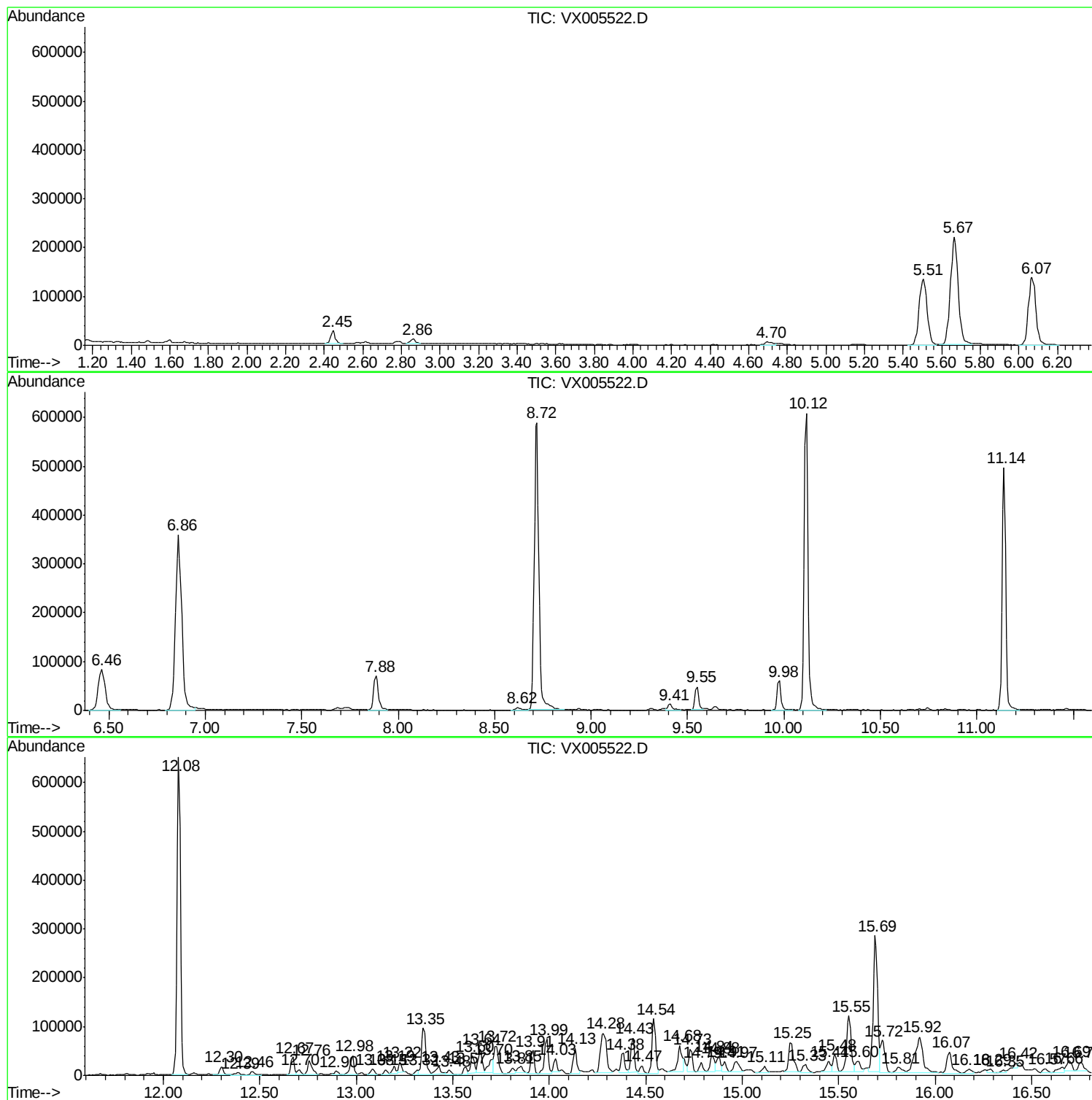
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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



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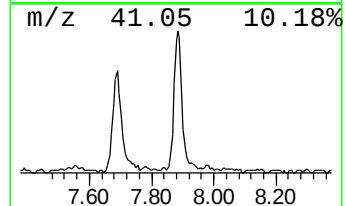
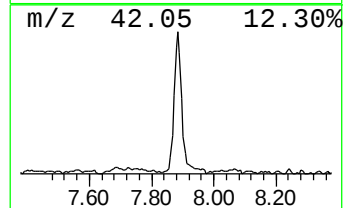
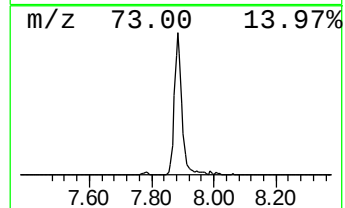
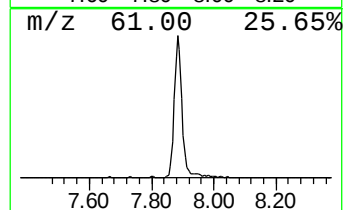
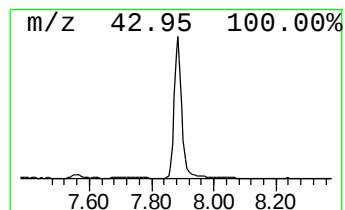
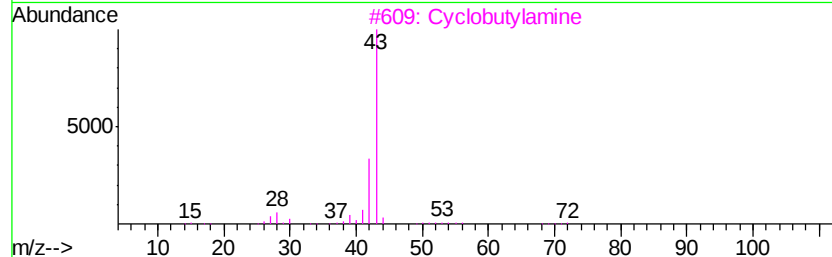
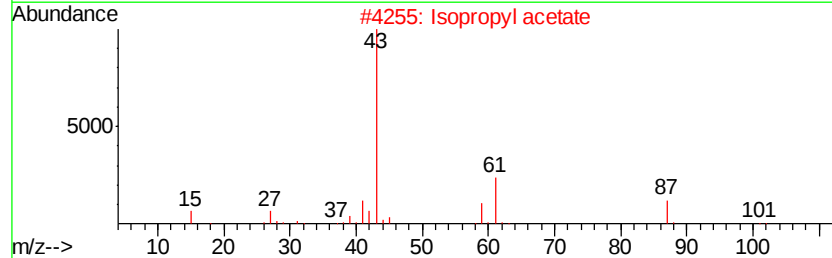
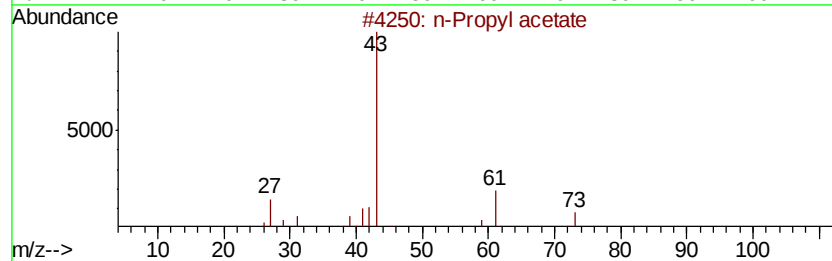
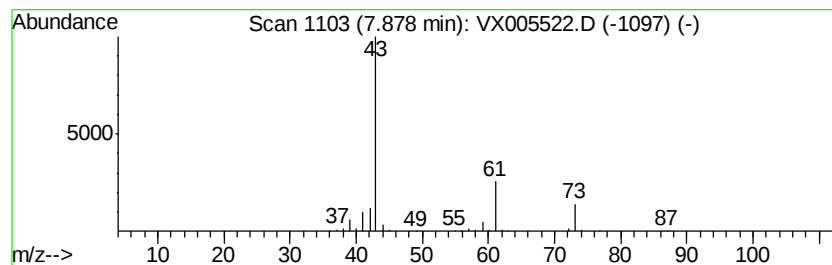
Instrument :
MSVOA_X
ClientSampled :
KG-V23608

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 1 n-Propyl acetate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.88	7.54 ug/l	123066	1,4-Difluorobenzene	6.86
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	n-Propyl acetate	102 C5H10O2	000109-60-4	78
2	Isopropyl acetate	102 C5H10O2	000108-21-4	33
3	Cyclobutylamine	71 C4H9N	002516-34-9	9
4	Acetic acid, pentyl ester	130 C7H14O2	000628-63-7	9
5	1-Butanamine	73 C4H11N	000109-73-9	7



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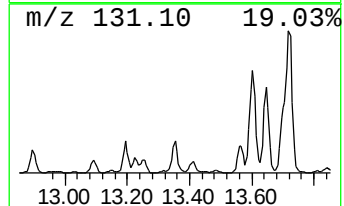
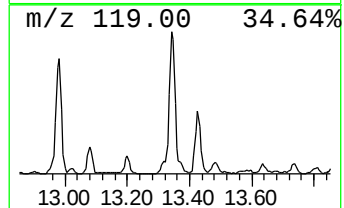
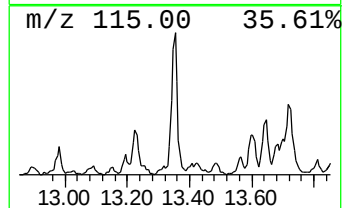
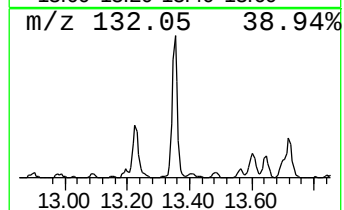
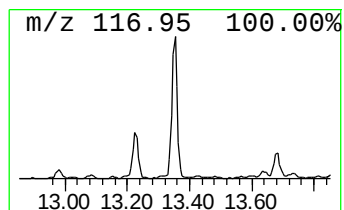
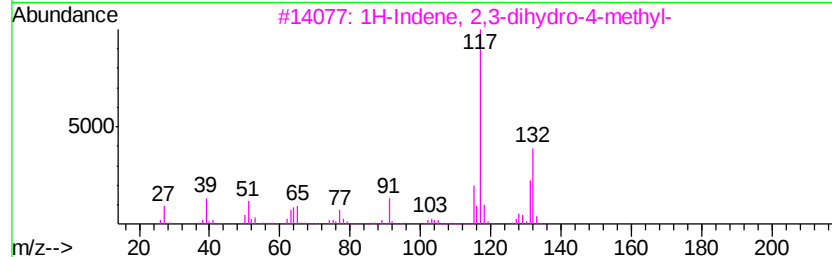
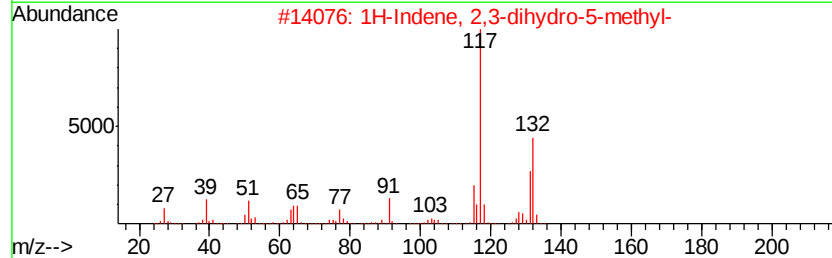
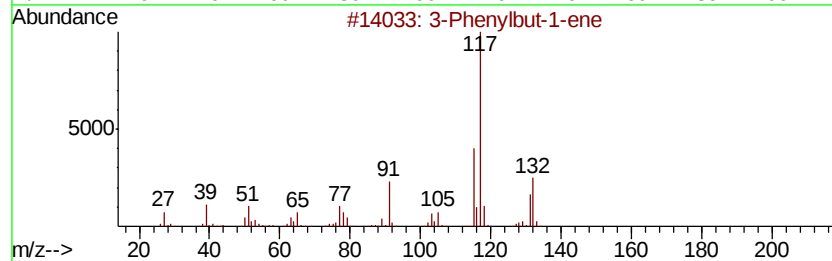
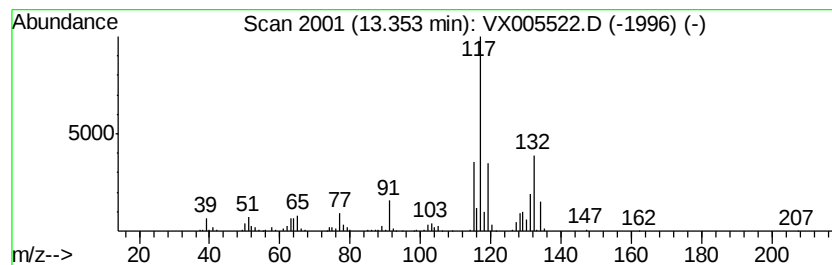
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Phenylbut-1-ene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.35	9.12 ug/l	145976	1,4-Dichlorobenzene-d4	12.08		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Phenylbut-1-ene	132	C10H12	000934-10-1	55
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	55
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	55
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	50
5		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	50



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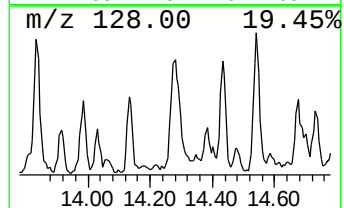
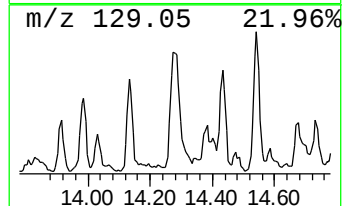
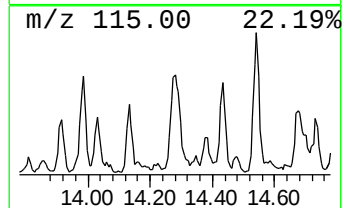
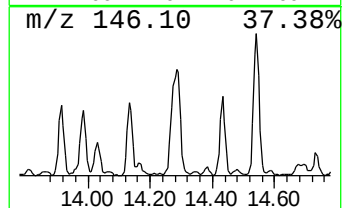
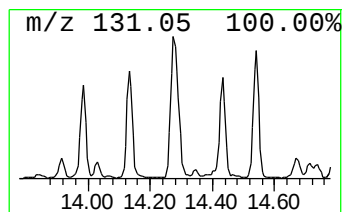
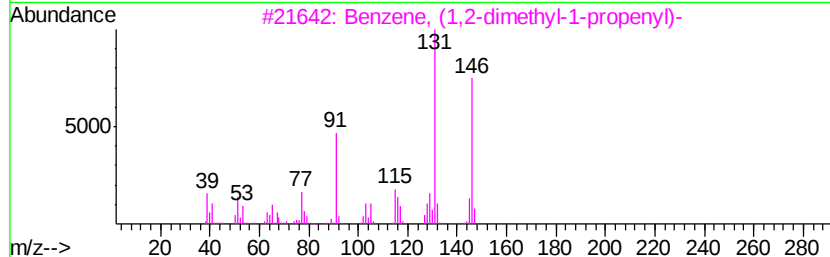
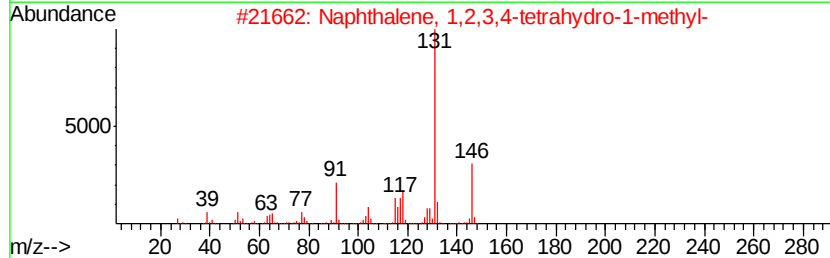
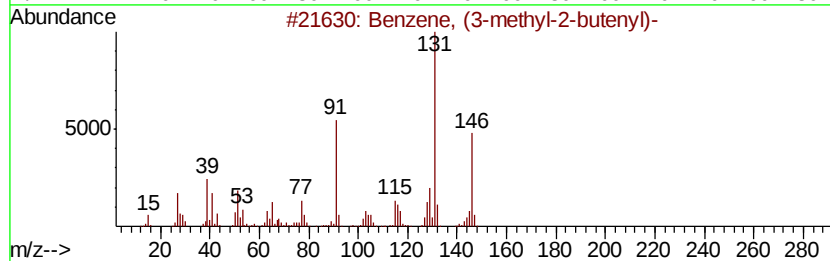
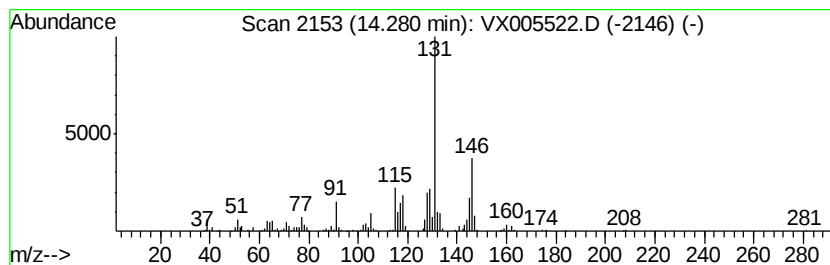
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, (3-methyl-2-butenyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	11.36 ug/l	181857	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	81
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76
3		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	76
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	76
5		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	74



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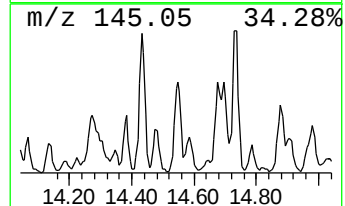
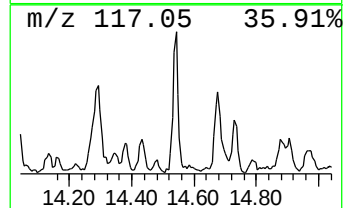
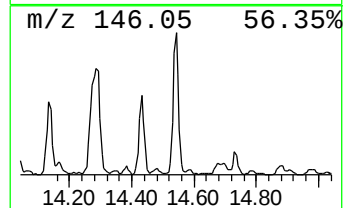
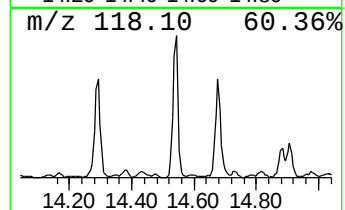
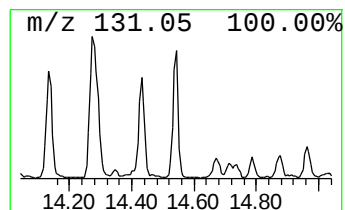
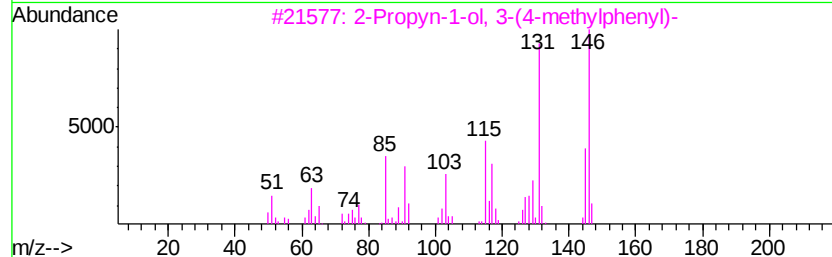
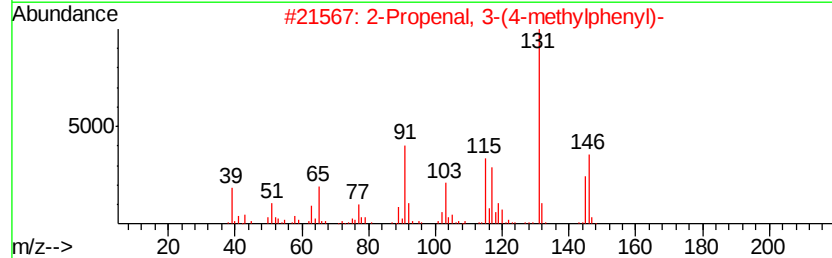
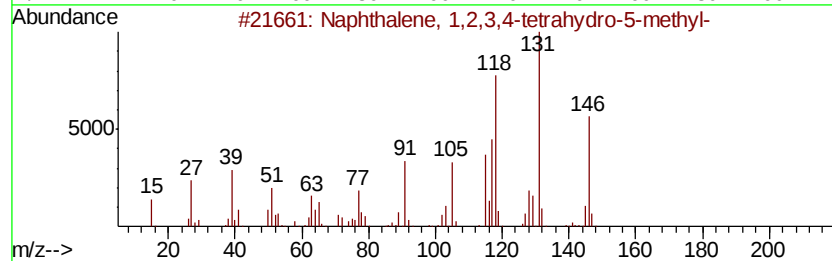
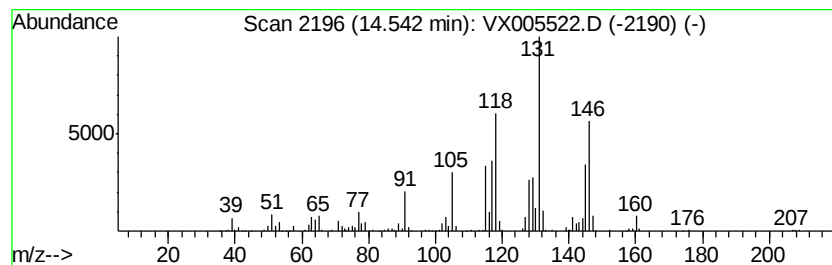
Instrument :
MSVOA_X
ClientSampleId :
KG-V23608

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.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 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	9.30 ug/l	148849	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 87
2		2-Propenal, 3-(4-methylphenyl)-	146 C10H10O	001504-75-2 64
3		2-Propyn-1-ol, 3-(4-methylphenyl)-	146 C10H10O	016017-24-6 64
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 64
5		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102318\
Data File : VX005522.D
Acq On : 23 Oct 2018 15:47
Operator : JC/MD
Sample : J5622-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
KG-V23608

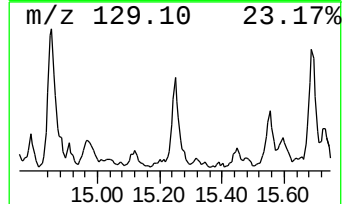
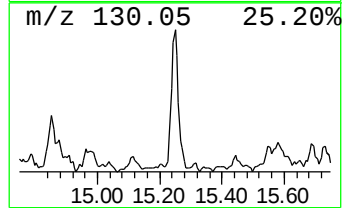
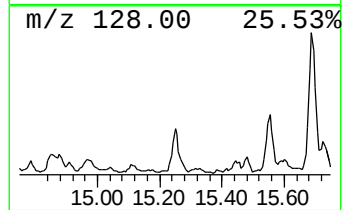
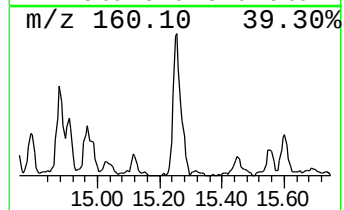
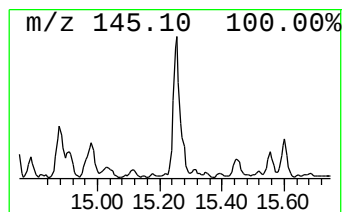
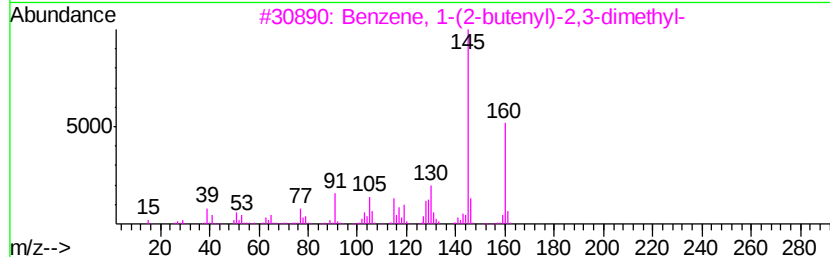
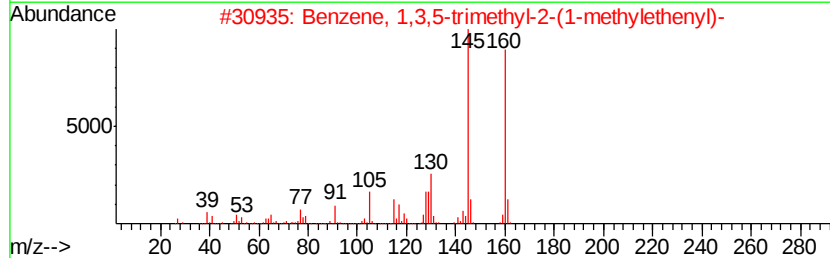
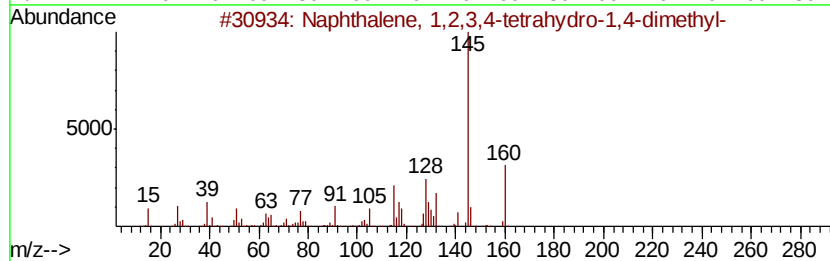
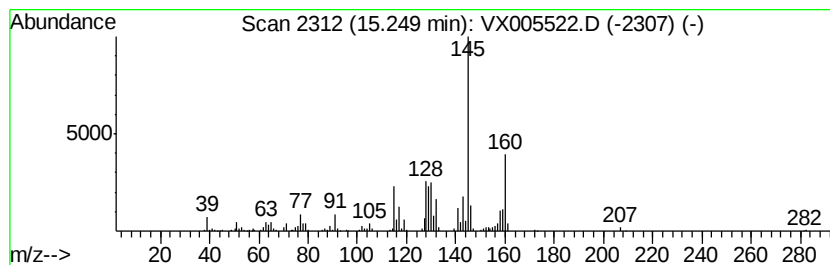
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	6.58 ug/l	105365	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	89
2		Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	81
3		Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	76
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	68
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	020027-77-4	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102318\
Data File : VX005522.D
Acq On : 23 Oct 2018 15:47
Operator : JC/MD
Sample : J5622-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KG-V23608

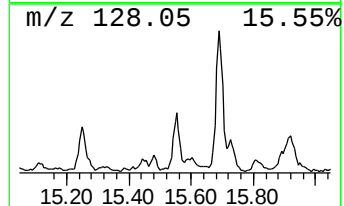
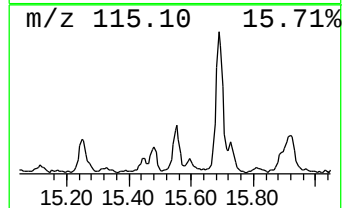
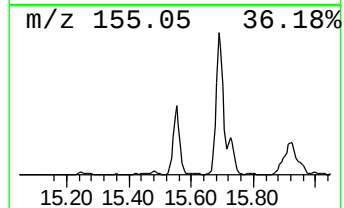
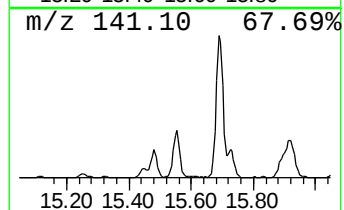
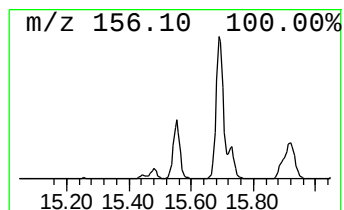
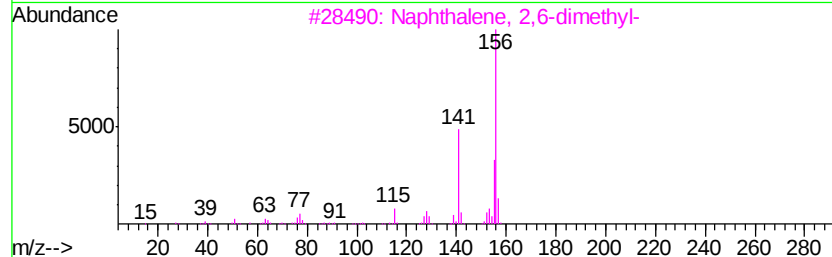
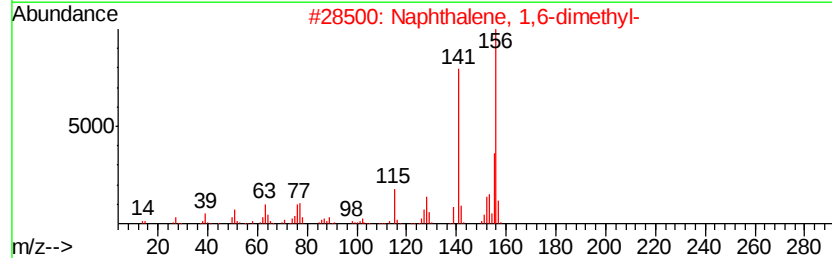
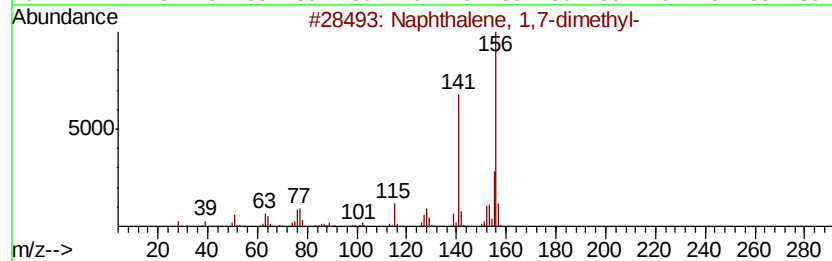
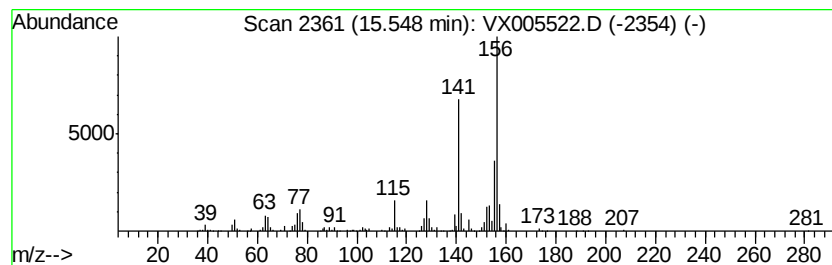
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,7-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	11.45 ug/l	183248	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102318\
Data File : VX005522.D
Acq On : 23 Oct 2018 15:47
Operator : JC/MD
Sample : J5622-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
KG-V23608

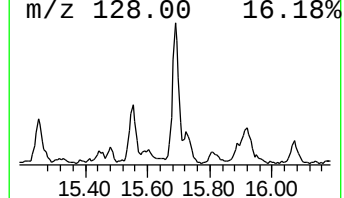
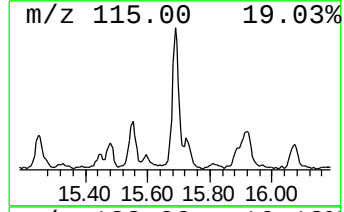
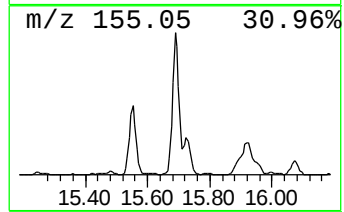
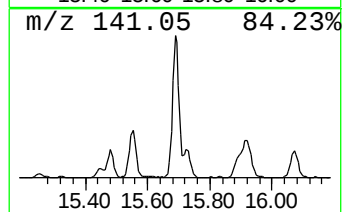
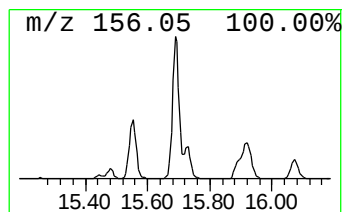
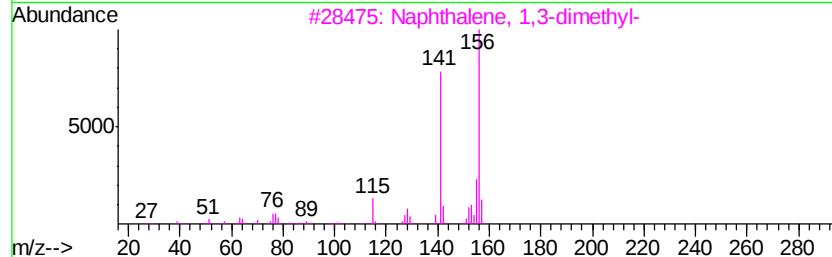
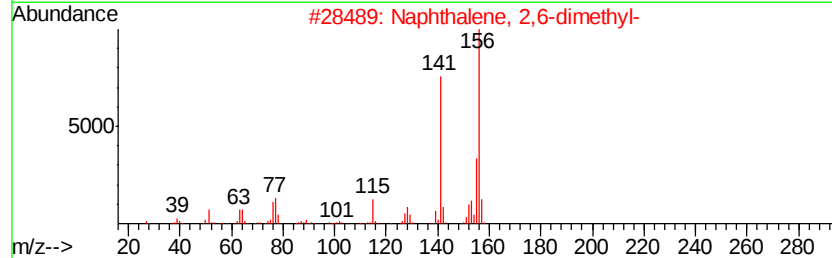
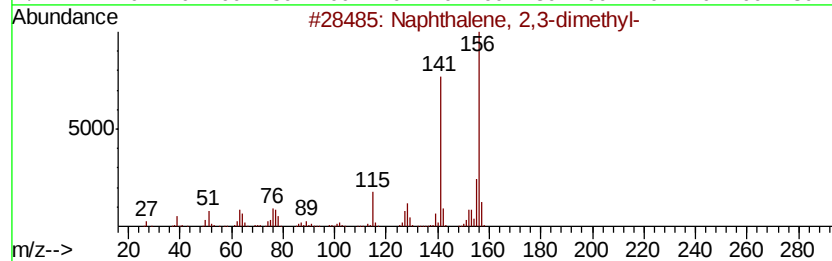
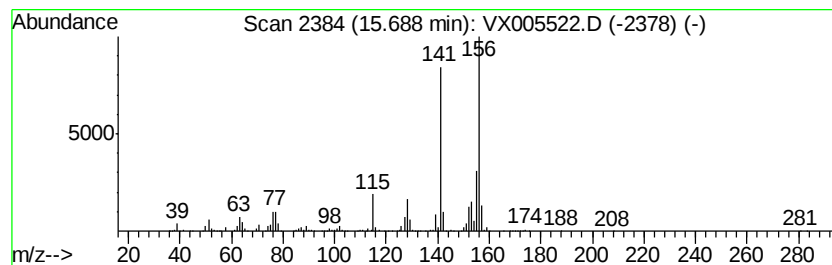
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	27.59 ug/l	441424	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	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 X\DATA\VX102318\
Data File : VX005522.D
Acq On : 23 Oct 2018 15:47
Operator : JC/MD
Sample : J5622-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

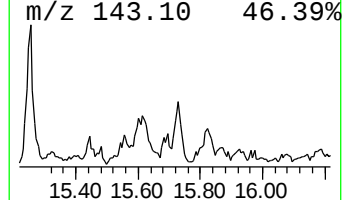
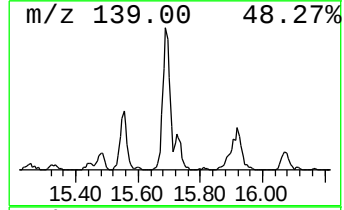
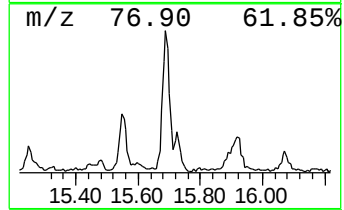
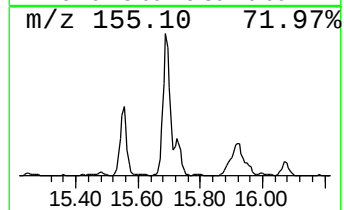
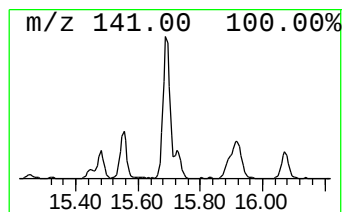
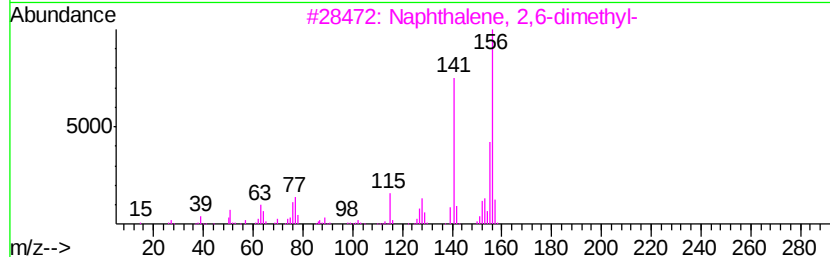
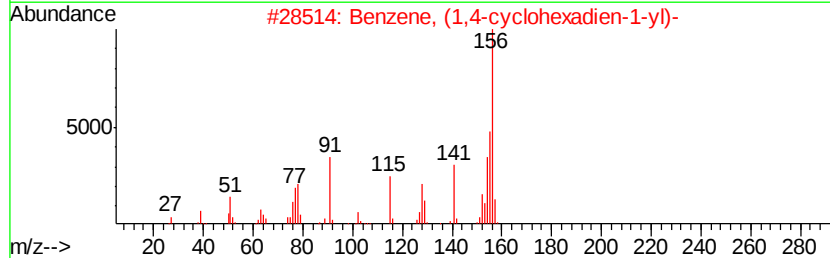
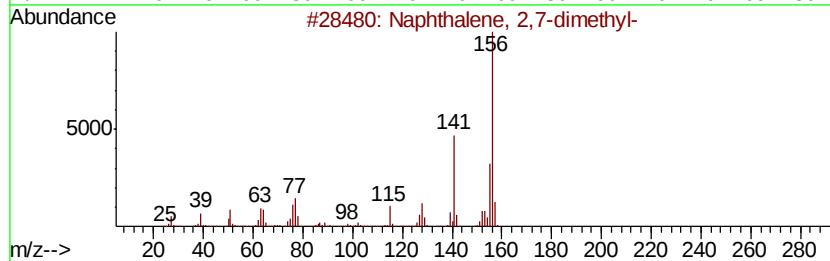
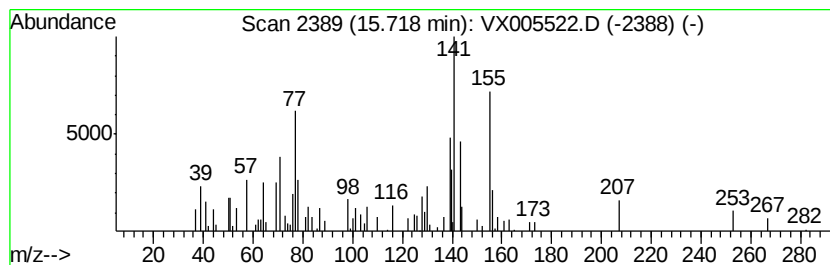
Instrument :
MSVOA_X
ClientSampleId :
KG-V23608

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	6.34 ug/l	101453	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 76
2		Benzene, (1,4-cyclohexadien-1-yl)-	156 C12H12	013703-52-1 70
3		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 68
4		Benzene, 2,5-cyclohexadien-1-yl-	156 C12H12	004794-05-2 58
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102318\
Data File : VX005522.D
Acq On : 23 Oct 2018 15:47
Operator : JC/MD
Sample : J5622-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

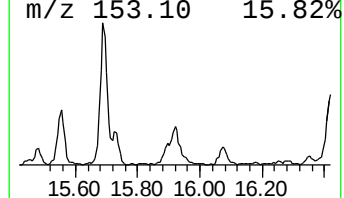
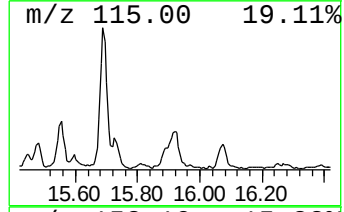
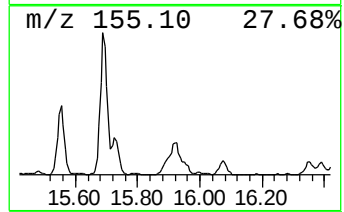
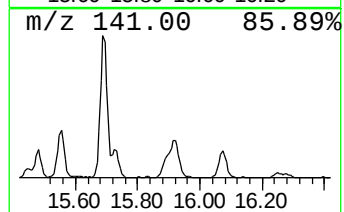
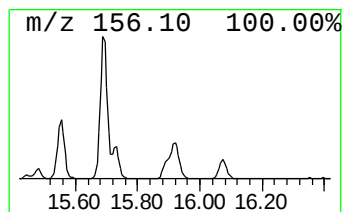
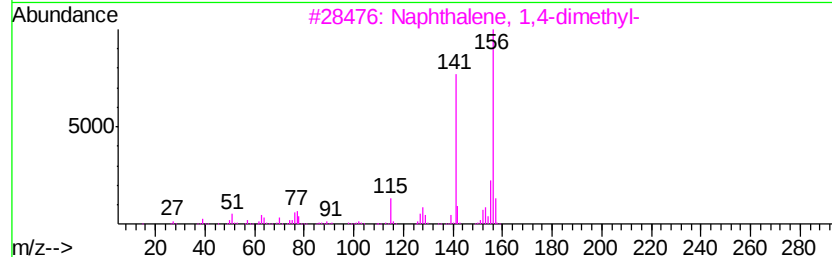
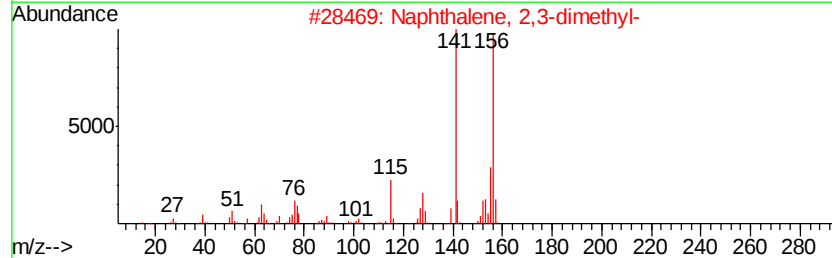
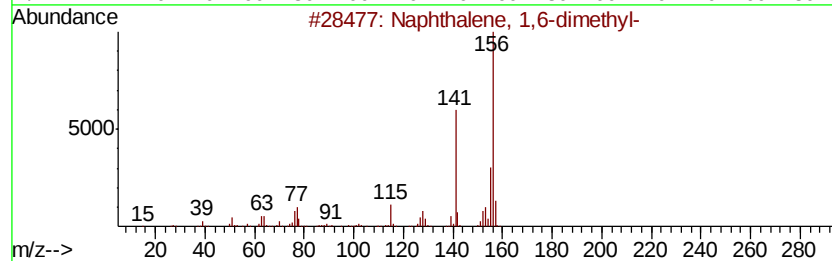
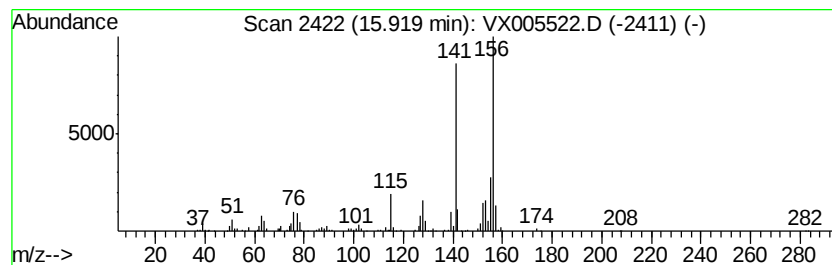
Instrument :
MSVOA_X
ClientSampleId :
KG-V23608

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.92	12.92 ug/l	206719	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX102318\
Data File : VX005522.D
Acq On : 23 Oct 2018 15:47
Operator : JC/MD
Sample : J5622-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KG-V23608

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.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
n-Propyl acetate	7.88	7.5	ug/l	123066	2	6.86	816594	50.0
3-Phenylbut-1-ene	13.35	9.1	ug/l	145976	4	12.08	800102	50.0
Benzene, (3-methy...	14.28	11.4	ug/l	181857	4	12.08	800102	50.0
Naphthalene, 1,2,...	14.54	9.3	ug/l	148849	4	12.08	800102	50.0
Naphthalene, 1,2,...	15.25	6.6	ug/l	105365	4	12.08	800102	50.0
Naphthalene, 1,7-...	15.55	11.4	ug/l	183248	4	12.08	800102	50.0
Naphthalene, 2,3-...	15.69	27.6	ug/l	441424	4	12.08	800102	50.0
Naphthalene, 2,7-...	15.72	6.3	ug/l	101453	4	12.08	800102	50.0
Naphthalene, 1,6-...	15.92	12.9	ug/l	206719	4	12.08	800102	50.0