

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082218\
 Data File : VN050821.D
 Acq On : 22 Aug 2018 15:51
 Operator : MD\SY
 Sample : J4586-01 10X
 Misc : 7.01µ/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EP1-275-UST-SLUDGE

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_N\METHODS\82N082118W.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	1.658	3	21	44	rBV	867191	1436022	23.25%	1.457%
2	7.593	1847	1867	1878	rBV	554829	1377820	22.31%	1.398%
3	7.667	1879	1890	1921	rVB	797787	1898504	30.74%	1.927%
4	8.030	1982	2003	2031	rBV	531402	1251208	20.26%	1.270%
5	8.262	2041	2075	2076	rBV3	63747	269266	4.36%	0.273%
6	8.587	2160	2176	2204	rBV	1210002	2499783	40.48%	2.537%
7	10.091	2629	2644	2678	rBV	2137290	3970382	64.29%	4.029%
8	11.204	2973	2990	3008	rVB6	50928	144292	2.34%	0.146%
9	11.300	3008	3020	3029	rBV2	48743	86933	1.41%	0.088%
10	11.410	3039	3054	3074	rBV	1574279	2802734	45.38%	2.844%
11	11.631	3102	3123	3138	rBV6	167932	464914	7.53%	0.472%
12	11.947	3200	3221	3237	rBV3	229849	625694	10.13%	0.635%
13	12.046	3243	3252	3260	rBV	198471	282076	4.57%	0.286%
14	12.120	3266	3275	3280	rBV3	94318	161294	2.61%	0.164%
15	12.159	3281	3287	3297	rVB4	137676	249904	4.05%	0.254%
16	12.300	3320	3331	3335	rBV6	81298	166087	2.69%	0.169%
17	12.339	3337	3343	3347	rVV	107547	137715	2.23%	0.140%
18	12.400	3354	3362	3372	rVV	1330143	2152088	34.85%	2.184%
19	12.445	3372	3376	3385	rVB	170772	209988	3.40%	0.213%
20	12.561	3405	3412	3427	rVB6	126079	246422	3.99%	0.250%
21	12.654	3427	3441	3449	rBV	812508	1460173	23.64%	1.482%
22	12.728	3455	3464	3476	rVB3	2473294	4024872	65.17%	4.085%
23	12.892	3502	3515	3525	rBV	1035901	1899369	30.76%	1.928%
24	12.963	3530	3537	3549	rVB	688133	1060089	17.17%	1.076%
25	13.040	3550	3561	3570	rBV	947041	1518397	24.59%	1.541%
26	13.091	3570	3577	3586	rVB3	156795	230122	3.73%	0.234%
27	13.146	3587	3594	3608	rVB6	163463	356621	5.77%	0.362%
28	13.236	3609	3622	3630	rBV2	866255	1546507	25.04%	1.569%
29	13.284	3631	3637	3644	rBV4	335414	458882	7.43%	0.466%
30	13.345	3644	3656	3661	rBV	1648437	2914906	47.20%	2.958%
31	13.426	3675	3681	3693	rVB2	225501	369223	5.98%	0.375%
32	13.512	3700	3708	3710	rBV	451531	535358	8.67%	0.543%
33	13.541	3711	3717	3723	rVV	1251757	1871750	30.31%	1.900%
34	13.583	3723	3730	3743	rVB2	1627670	2672434	43.27%	2.712%

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35	13.673	3746	3758	3767	rBV	1834398	2911108	47.14%	2.954%
36	13.741	3771	3779	3788	rVB	835567	1223420	19.81%	1.242%
37	13.802	3790	3798	3802	rBV	709573	1042366	16.88%	1.058%
38	13.831	3803	3807	3816	rVB	1010107	1395172	22.59%	1.416%
39	13.889	3817	3825	3834	rVB	1598676	2401617	38.89%	2.437%
40	13.943	3835	3842	3848	rBV2	255255	379065	6.14%	0.385%
41	13.995	3848	3858	3868	rVB	1839926	2951168	47.79%	2.995%
42	14.053	3868	3876	3878	rBV	229998	310581	5.03%	0.315%
43	14.127	3890	3899	3907	rBV2	732839	1237861	20.04%	1.256%
44	14.210	3919	3925	3930	rBV	486691	621355	10.06%	0.631%
45	14.249	3931	3937	3945	rVB	1497068	2118810	34.31%	2.150%
46	14.297	3945	3952	3958	rBV2	693112	931848	15.09%	0.946%
47	14.332	3958	3963	3971	rVB2	476979	587860	9.52%	0.597%
48	14.432	3986	3994	4001	rVB3	529140	827703	13.40%	0.840%
49	14.477	4001	4008	4016	rBV3	558747	883999	14.31%	0.897%
50	14.525	4016	4023	4031	rVB	976925	1367539	22.14%	1.388%
51	14.586	4031	4042	4055	rBV5	2763607	6175544	100.00%	6.267%
52	14.647	4056	4061	4072	rVB2	540123	858238	13.90%	0.871%
53	14.741	4082	4090	4100	rVB	1228382	1826261	29.57%	1.853%
54	14.853	4107	4125	4131	rBV3	1243385	2899902	46.96%	2.943%
55	14.959	4148	4158	4174	rVB2	1276346	2696104	43.66%	2.736%
56	15.110	4194	4205	4212	rBV5	274847	642886	10.41%	0.652%
57	15.191	4221	4230	4238	rBV	973378	1576650	25.53%	1.600%
58	15.249	4239	4248	4268	rVB7	734362	2568927	41.60%	2.607%
59	15.419	4289	4301	4314	rBV3	834206	1792794	29.03%	1.819%
60	15.577	4335	4350	4353	rBV5	700870	1486811	24.08%	1.509%
61	15.615	4355	4362	4377	rVB	1704148	3063953	49.61%	3.109%
62	15.699	4382	4388	4398	rVV4	462034	910196	14.74%	0.924%
63	15.770	4400	4410	4435	rVB6	700647	2156192	34.92%	2.188%
64	15.914	4443	4455	4466	rBV2	1197253	2388887	38.68%	2.424%
65	16.101	4504	4513	4524	rVB3	894568	1698987	27.51%	1.724%
66	16.162	4525	4532	4542	rVB3	532904	890068	14.41%	0.903%
67	16.355	4579	4592	4604	rBV4	951097	2361770	38.24%	2.397%

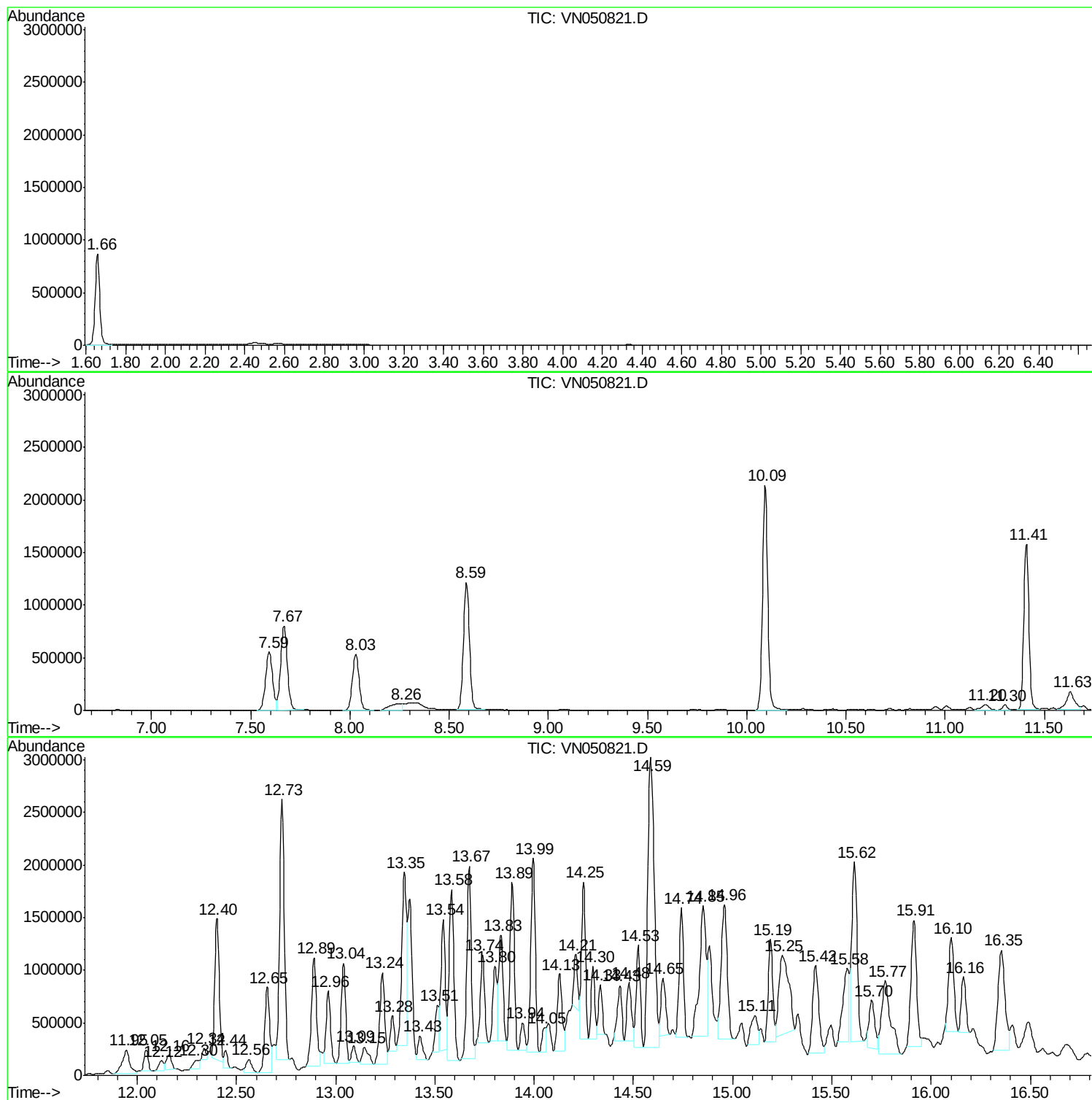
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Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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



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ALS Vial : 14 Sample Multiplier: 1

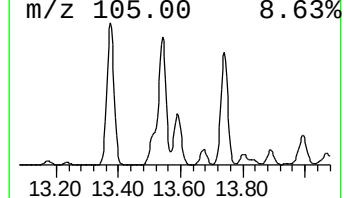
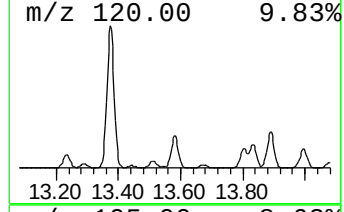
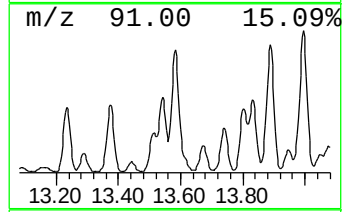
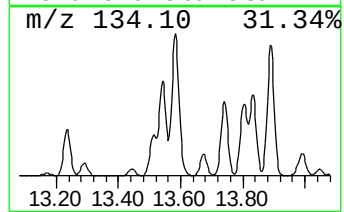
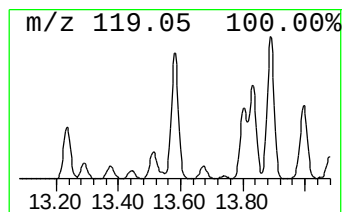
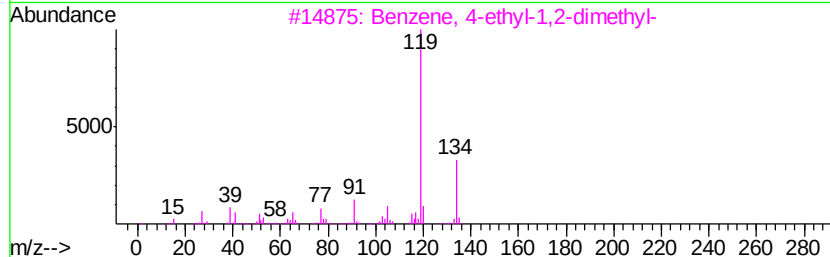
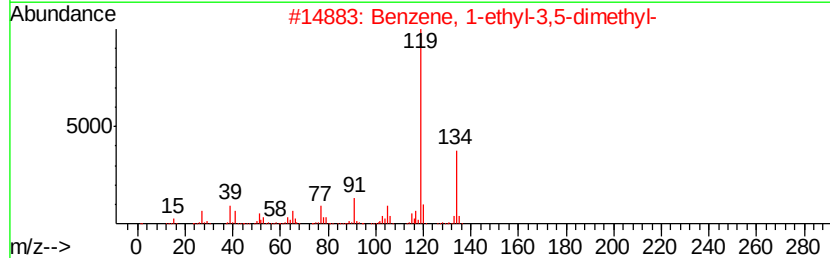
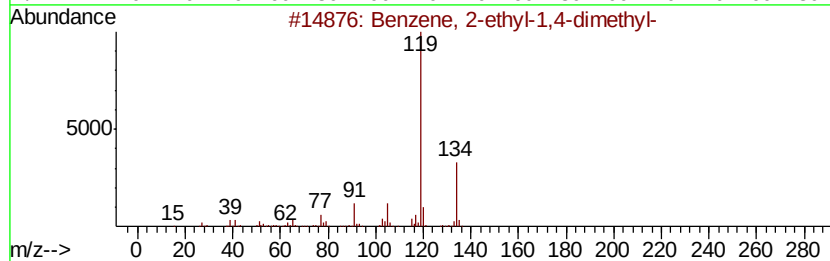
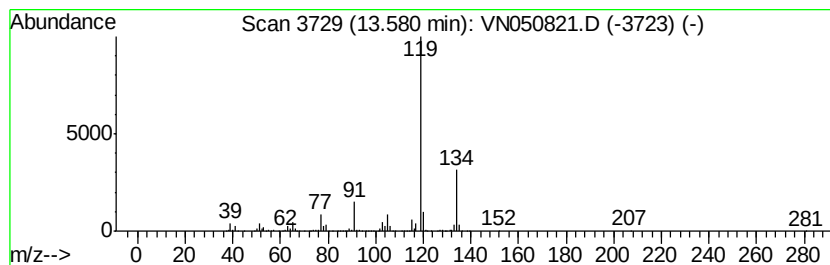
Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.58	45.84 ug/l	2672430	1,4-Dichlorobenzene-d4	13.35		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



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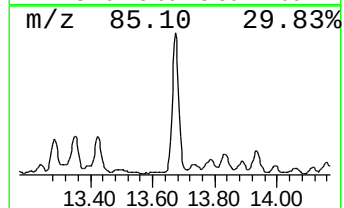
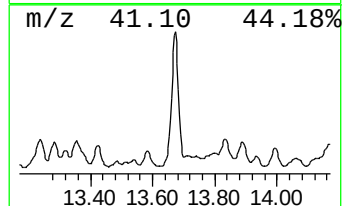
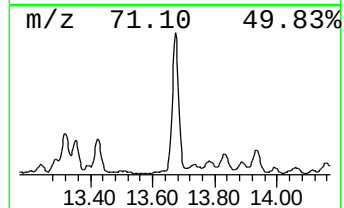
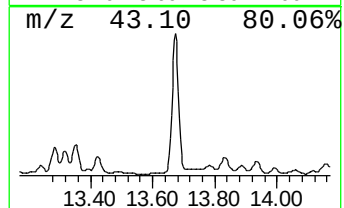
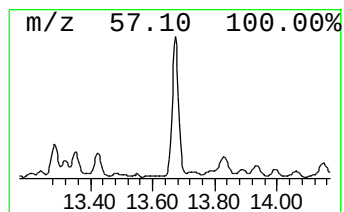
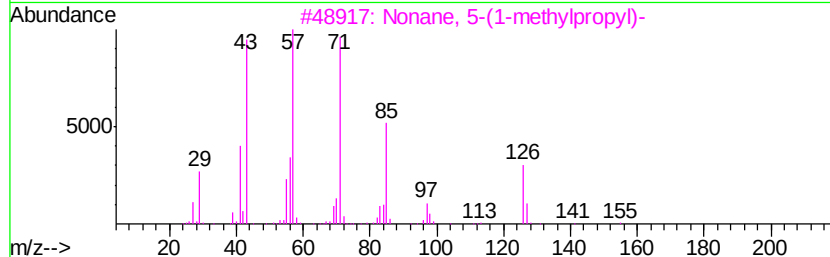
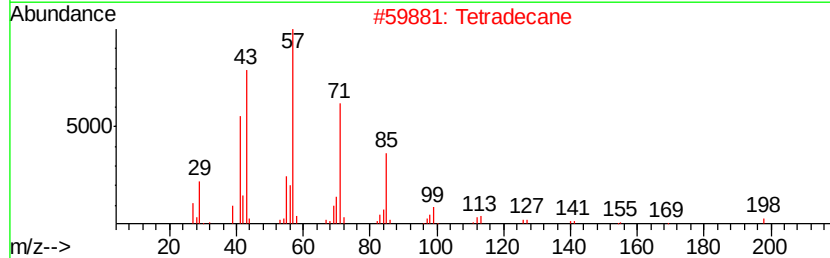
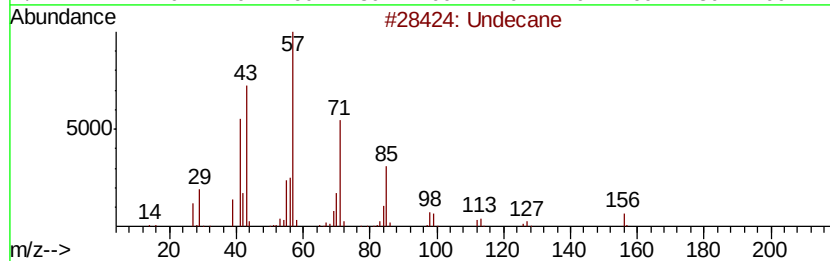
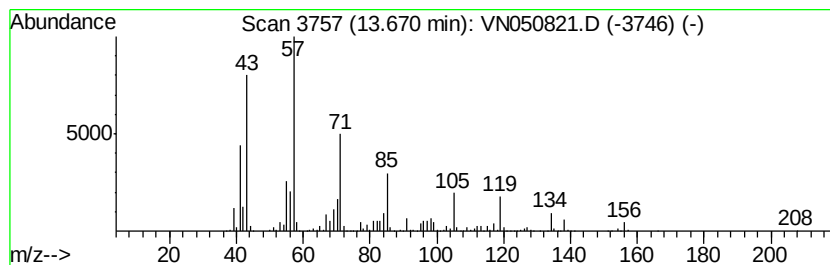
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Undecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	49.93 ug/l	2911110	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane	156 C11H24	001120-21-4	92
2	Tetradecane	198 C14H30	000629-59-4	52
3	Nonane, 5-(1-methylpropyl)-	184 C13H28	062185-54-0	47
4	Hexadecane	226 C16H34	000544-76-3	47
5	Tridecane	184 C13H28	000629-50-5	47



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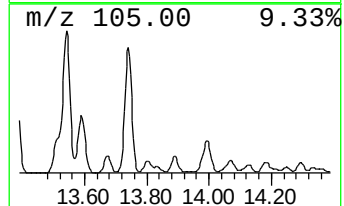
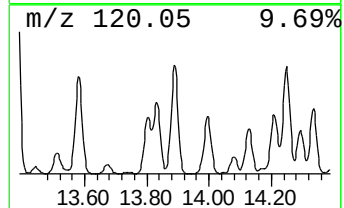
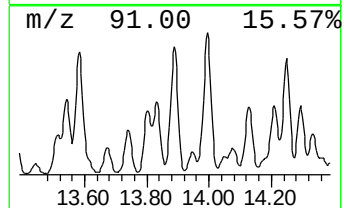
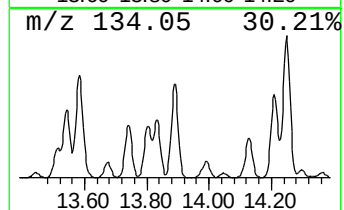
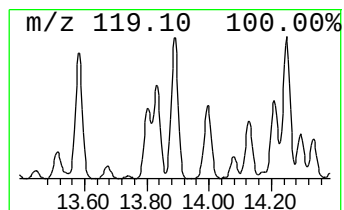
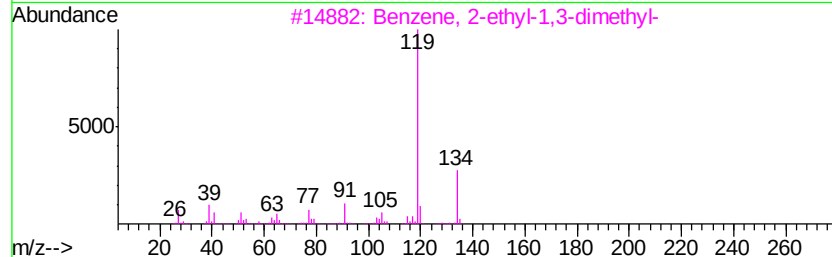
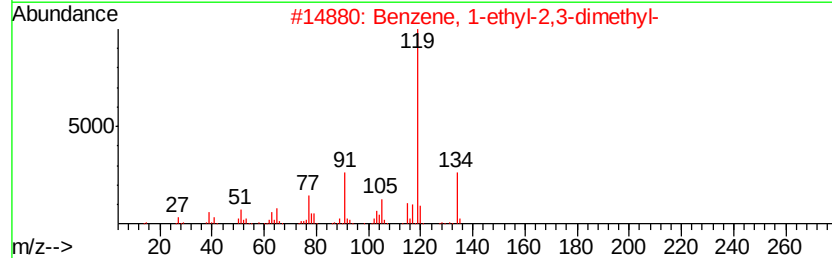
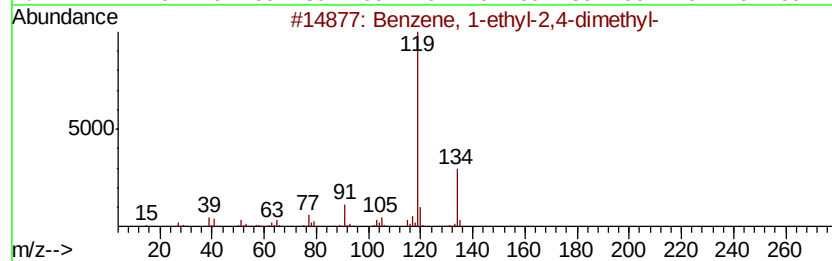
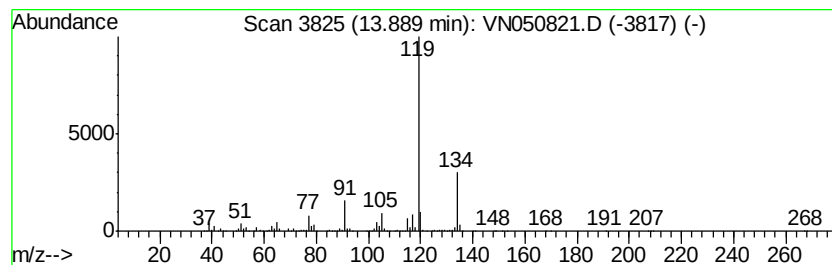
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.89	41.20 ug/l	2401620	1,4-Dichlorobenzene-d4	13.35	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



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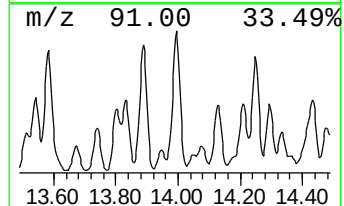
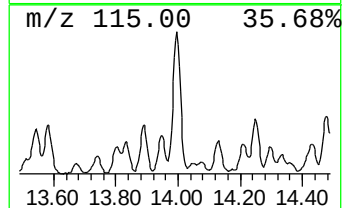
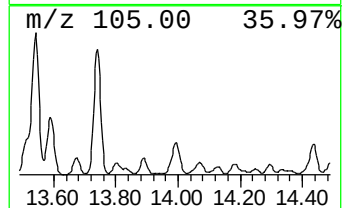
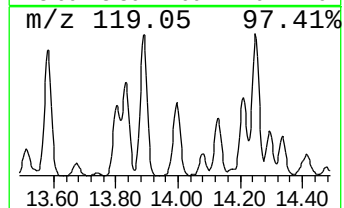
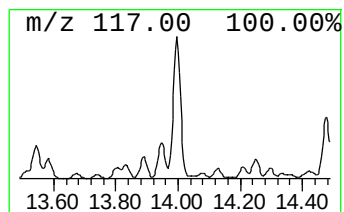
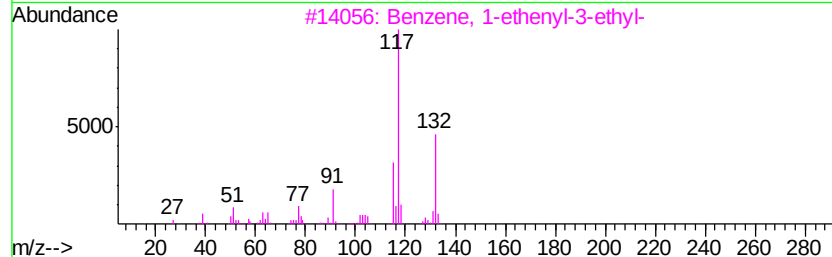
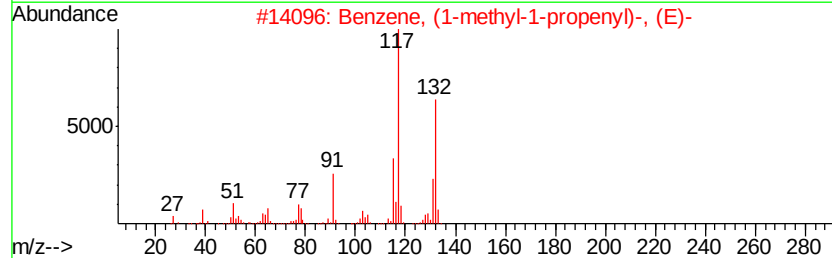
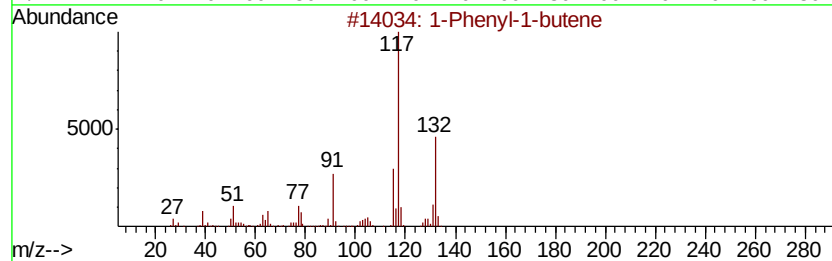
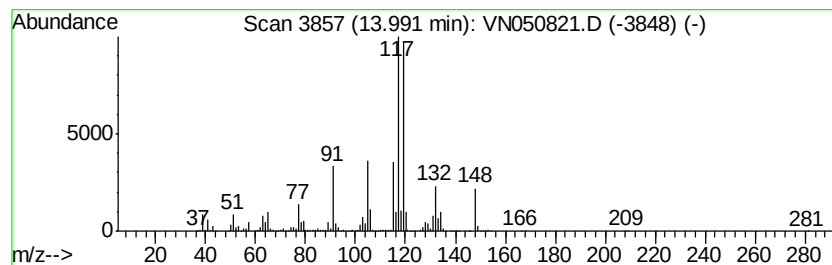
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Peak Number 4 1-Phenyl-1-butene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	50.62 ug/l	2951170	1,4-Dichlorobenzene-d4	13.35

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	64
2	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	64
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	50
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
5	Indan, 1-methyl-	132	C10H12	000767-58-8	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082218\
Data File : VN050821.D
Acq On : 22 Aug 2018 15:51
Operator : MD\SY
Sample : J4586-01 10X
Misc : 7.01a/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

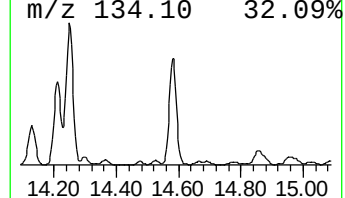
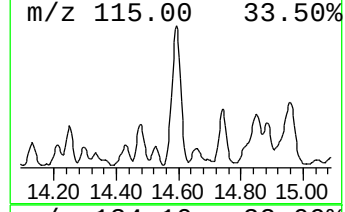
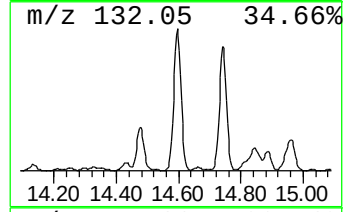
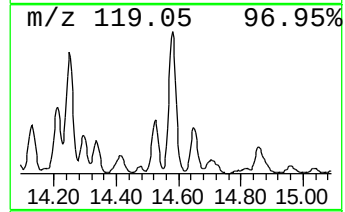
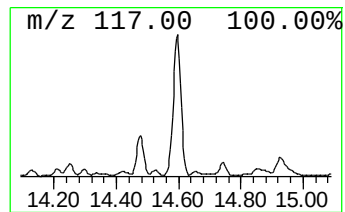
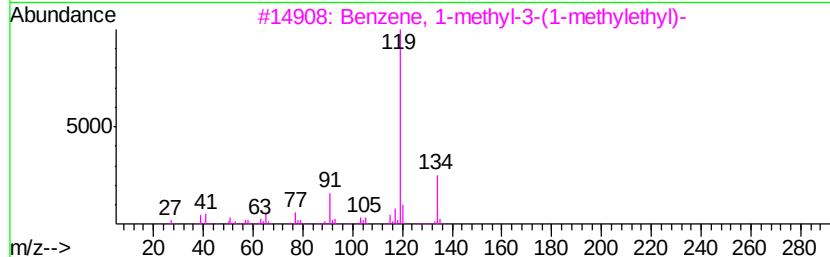
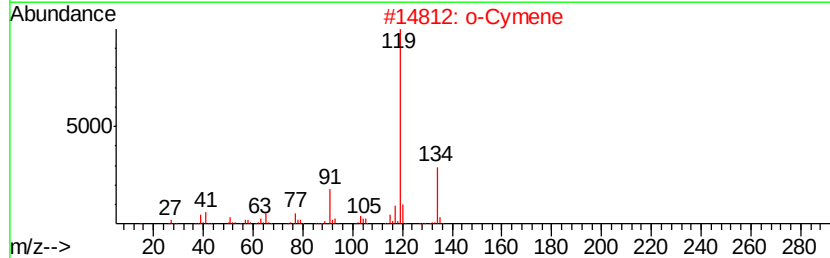
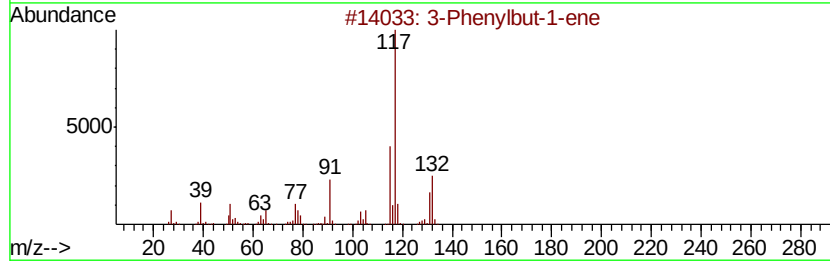
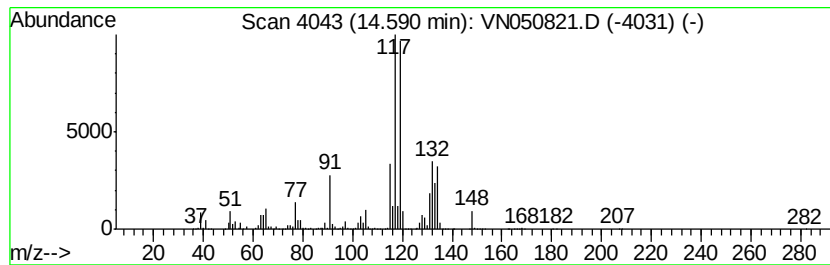
Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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

Peak Number 5 3-Phenylbut-1-ene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	105.93 ug/l	6175540	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Phenylbut-1-ene	132 C10H12	000934-10-1 80
2		o-Cymene	134 C10H14	000527-84-4 55
3		Benzene, 1-methyl-3-(1-methylethyl-)	134 C10H14	000535-77-3 55
4		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 55
5		p-Cymene	134 C10H14	000099-87-6 50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN082218\
Data File : VN050821.D
Acq On : 22 Aug 2018 15:51
Operator : MD\SY
Sample : J4586-01 10X
Misc : 7.01µ/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

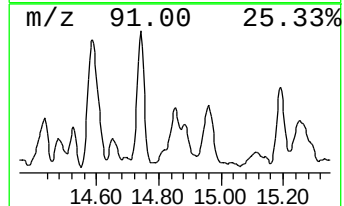
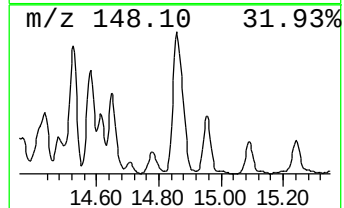
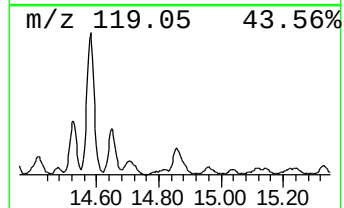
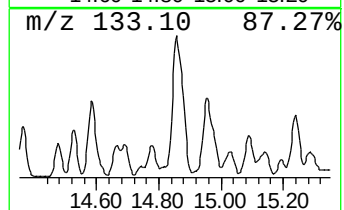
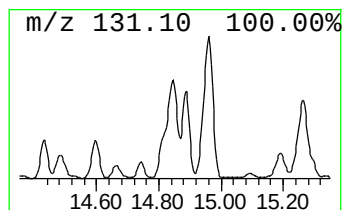
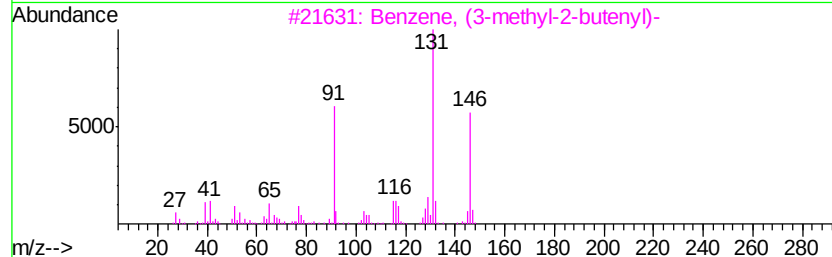
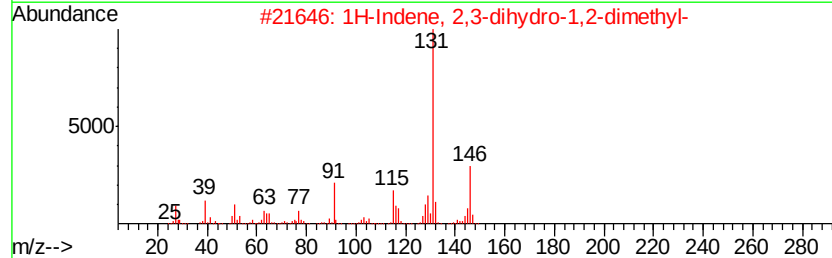
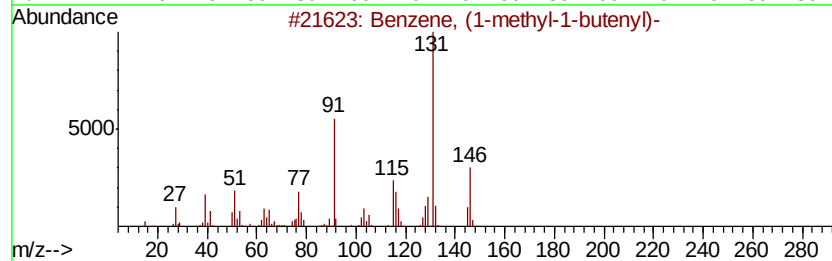
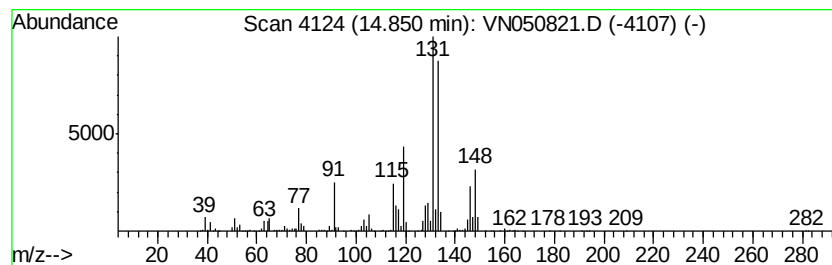
Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, (1-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	49.74 ug/l	2899900	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 87
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 62
3		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 59
4		Benzene, 2,4-diethyl-1-methyl-	148 C11H16	001758-85-6 55
5		Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1 50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN082218\
Data File : VN050821.D
Acq On : 22 Aug 2018 15:51
Operator : MD\SY
Sample : J4586-01 10X
Misc : 7.01µ/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

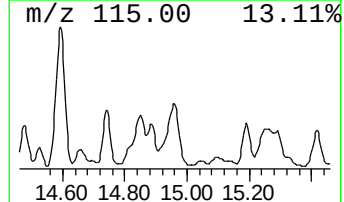
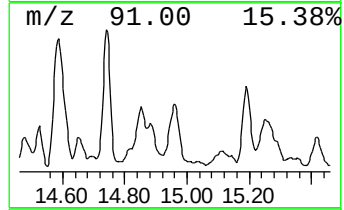
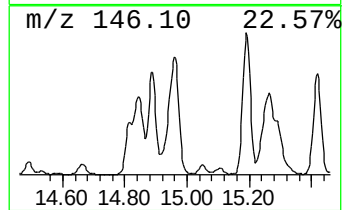
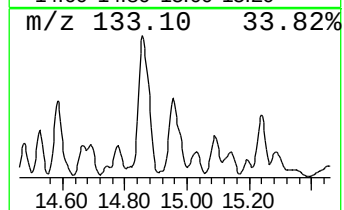
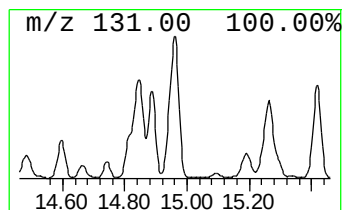
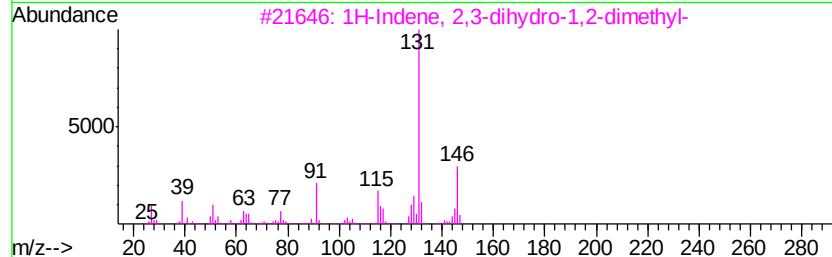
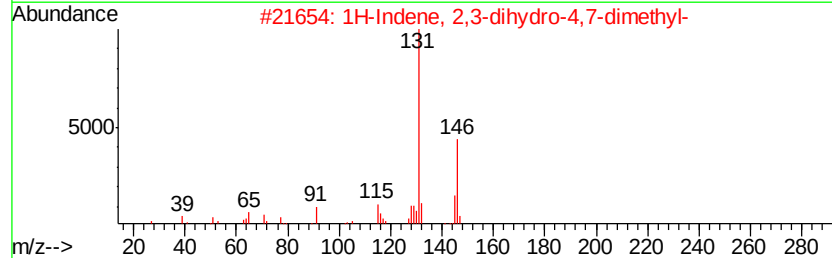
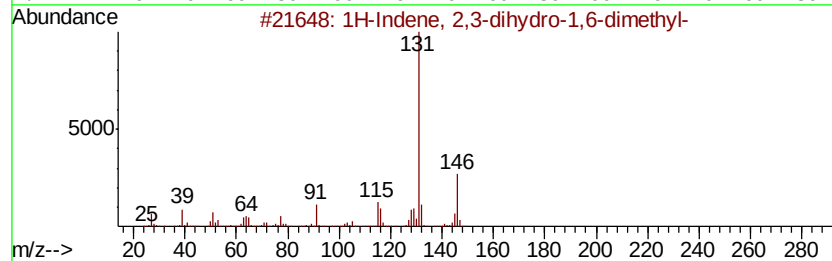
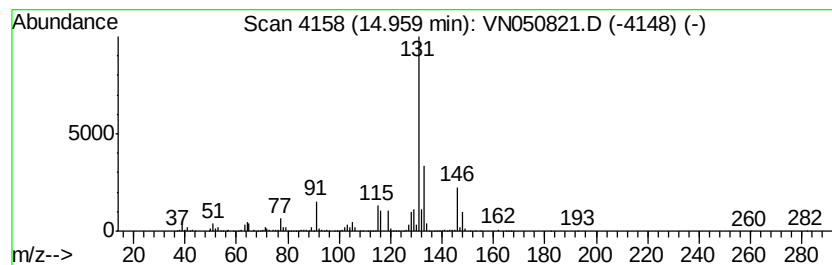
Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	46.25 ug/l	2696100	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	93
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	76
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8	76
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3	76
5	Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3	76



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN082218\
Data File : VN050821.D
Acq On : 22 Aug 2018 15:51
Operator : MD\SY
Sample : J4586-01 10X
Misc : 7.01µ/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

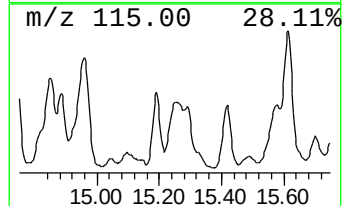
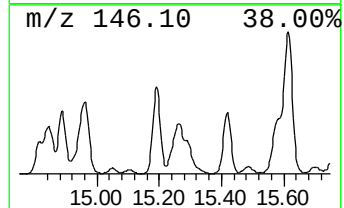
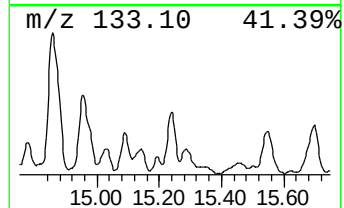
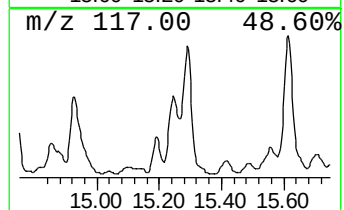
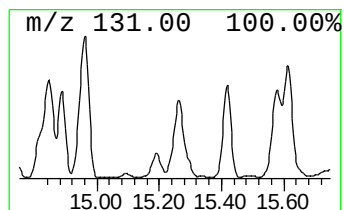
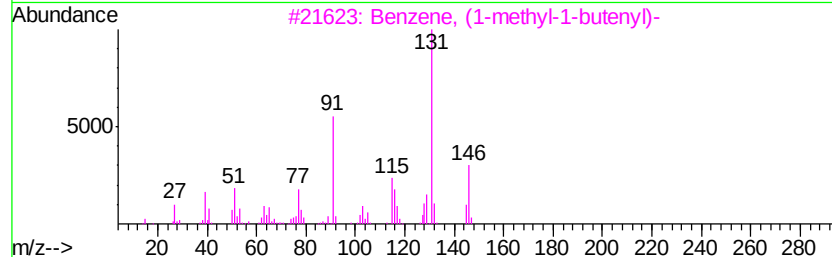
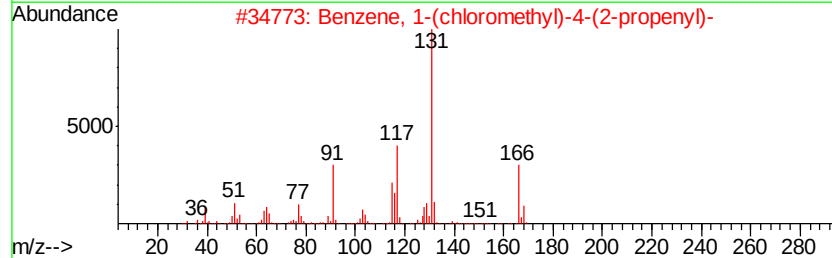
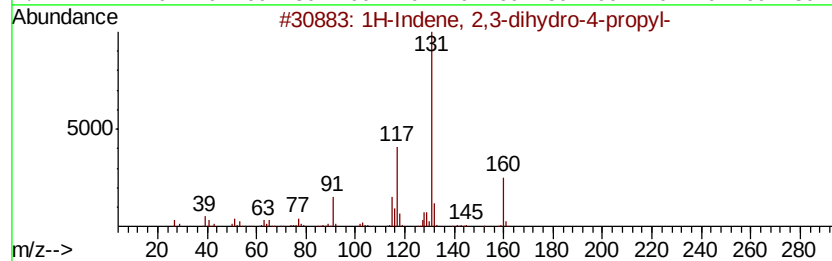
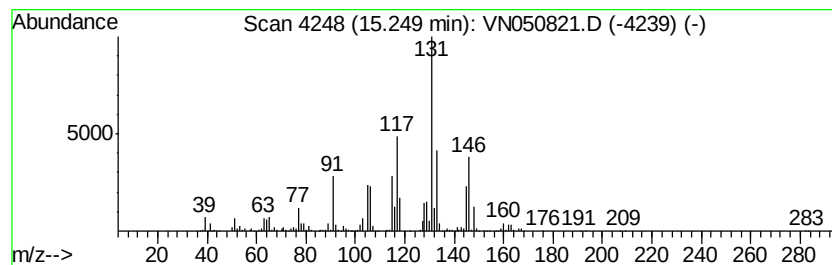
Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4-pr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	44.07 ug/l	2568930	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 2,3-dihydro-4-propyl-	160 C12H16	092013-16-6 87
2		Benzene, 1-(chloromethyl)-4-(2-p...	166 C10H11Cl	036875-10-2 55
3		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 55
4		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 55
5		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 55



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN082218\
Data File : VN050821.D
Acq On : 22 Aug 2018 15:51
Operator : MD\SY
Sample : J4586-01 10X
Misc : 7.01µ/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

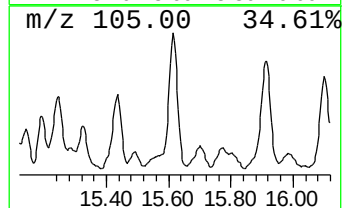
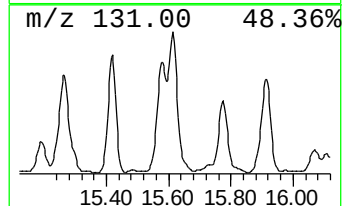
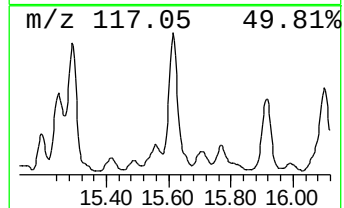
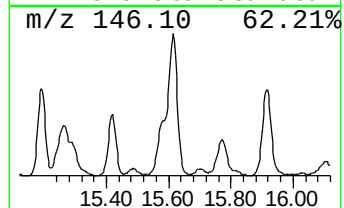
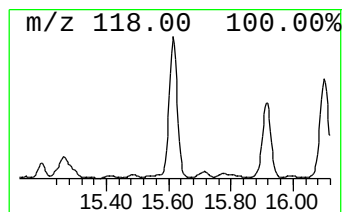
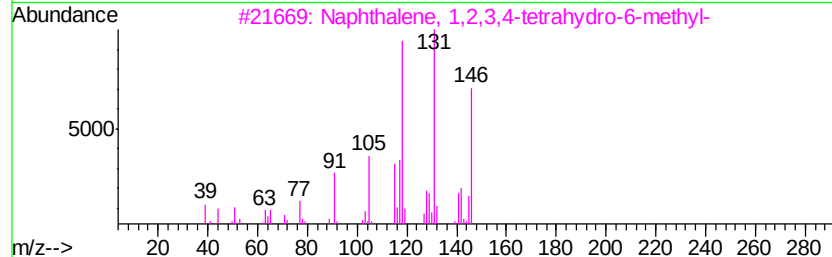
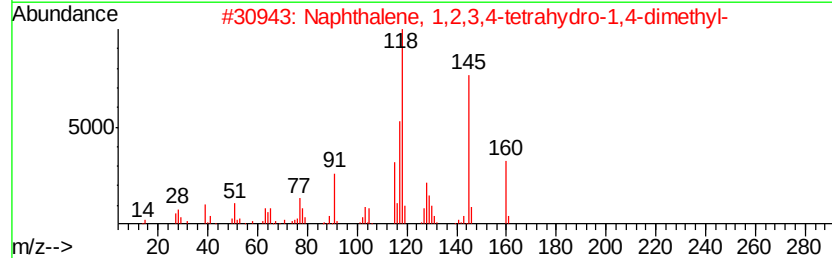
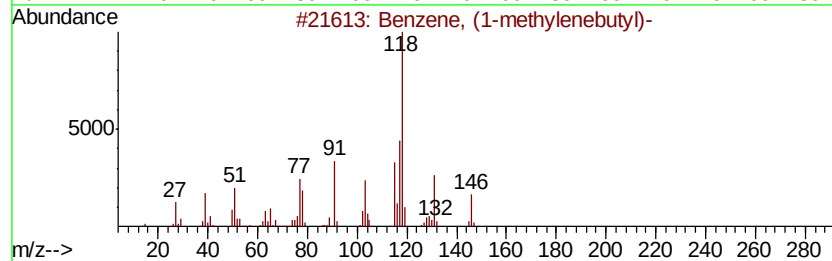
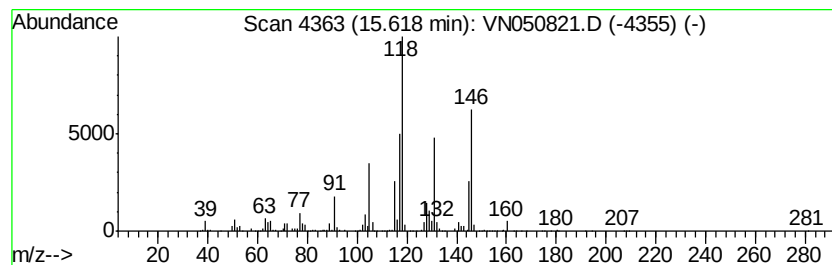
Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1-methylenebutyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.62	52.56 ug/l	3063950	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methylenebutyl)-	146 C11H14	005676-32-4 80
2		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	004175-54-6 64
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 58
4		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	013065-07-1 52
5		5H-Benzocycloheptene,6,7,8,9-tet...	146 C11H14	001075-16-7 43



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN082218\
Data File : VN050821.D
Acq On : 22 Aug 2018 15:51
Operator : MD\SY
Sample : J4586-01 10X
Misc : 7.01µ/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

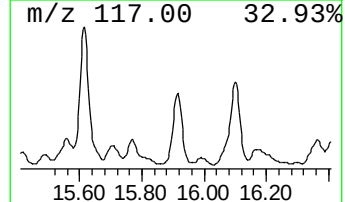
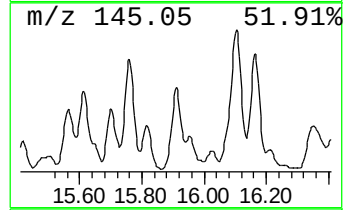
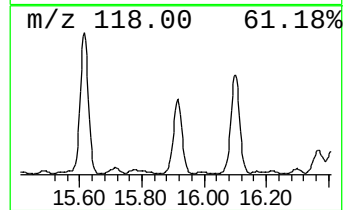
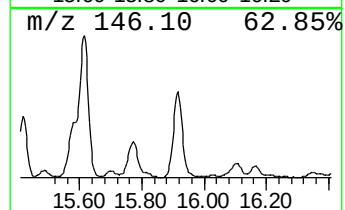
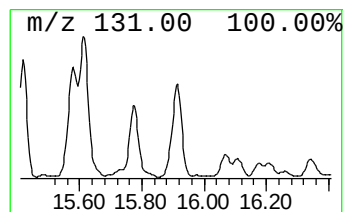
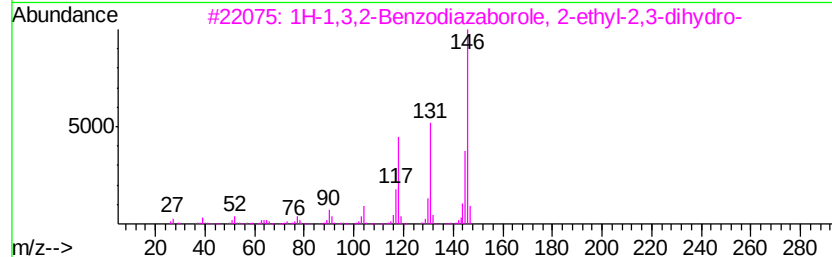
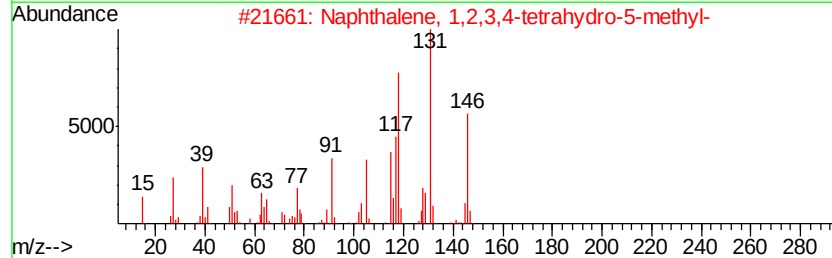
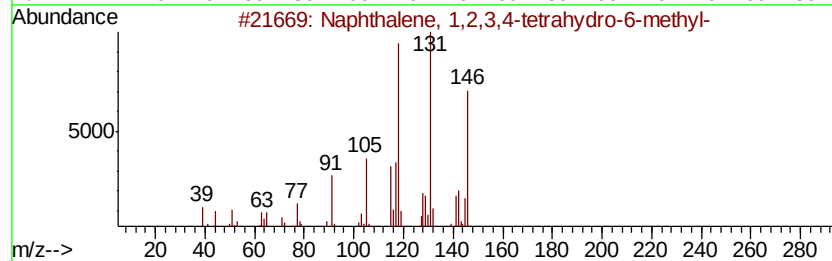
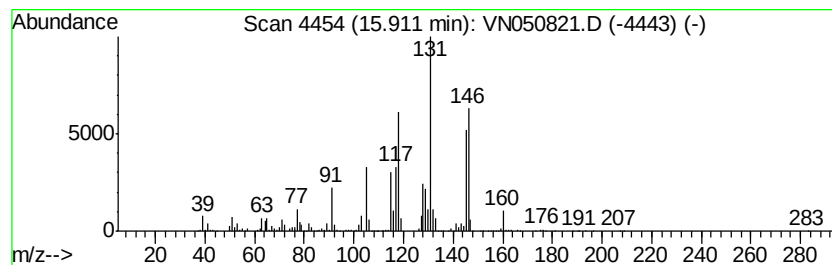
Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	40.98 ug/l	2388890	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 91
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 81
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 76
4		1H-Benzimidazole, 5,6-dimethyl-	146 C9H10N2	000582-60-5 64
5		2-Propyn-1-ol, 3-(4-methylphenyl)-	146 C10H10O	016017-24-6 62



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN082218\
Data File : VN050821.D
Acq On : 22 Aug 2018 15:51
Operator : MD\SY
Sample : J4586-01 10X
Misc : 7.01g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
EP1-275-UST-SLUDGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.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, 2-ethyl-...	13.58	45.8	ug/l	2672430	4	13.35	2914910	50.0
Undecane	13.67	49.9	ug/l	2911110	4	13.35	2914910	50.0
Benzene, 1-ethyl-...	13.89	41.2	ug/l	2401620	4	13.35	2914910	50.0
1-Phenyl-1-butene	13.99	50.6	ug/l	2951170	4	13.35	2914910	50.0
3-Phenylbut-1-ene	14.59	105.9	ug/l	6175540	4	13.35	2914910	50.0
Benzene, (1-methy...	14.85	49.7	ug/l	2899900	4	13.35	2914910	50.0
1H-Indene, 2,3-di...	14.96	46.3	ug/l	2696100	4	13.35	2914910	50.0
1H-Indene, 2,3-di...	15.25	44.1	ug/l	2568930	4	13.35	2914910	50.0
Benzene, (1-methy...	15.62	52.6	ug/l	3063950	4	13.35	2914910	50.0
Naphthalene, 1,2,...	15.91	41.0	ug/l	2388890	4	13.35	2914910	50.0