

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
 Data File : BM008829.D
 Acq On : 24 Jan 2017 07:44
 Operator : UM/SJ
 Sample : I1323-22
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA9R7

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 1 % 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:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.440	648	655	663	rVB	6460711	10901048	1.71%	0.554%
2	7.081	760	764	772	rVB2	5294602	9062711	1.42%	0.461%
3	7.322	790	805	808	rBV	6017522	13181724	2.07%	0.670%
4	7.369	808	813	820	rVB	14500558	22240733	3.49%	1.131%
5	8.086	914	935	939	rBV2	56905509	244657561	38.44%	12.444%
6	8.422	979	992	998	rBV	28568493	48805280	7.67%	2.482%
7	8.492	998	1004	1009	rVB	5449038	8301950	1.30%	0.422%
8	9.816	1222	1229	1232	rBV	5516013	10323226	1.62%	0.525%
9	10.845	1390	1404	1408	rVB2	55818197	146016189	22.94%	7.427%
10	10.933	1414	1419	1428	rVB	4102257	7600092	1.19%	0.387%
11	11.027	1428	1435	1441	rBV	4400495	9530296	1.50%	0.485%
12	11.104	1441	1448	1456	rVB	3684128	7873316	1.24%	0.400%
13	11.486	1506	1513	1524	rBV2	4883822	8943074	1.41%	0.455%
14	11.622	1531	1536	1540	rBV	9598332	14901768	2.34%	0.758%
15	11.663	1540	1543	1553	rVB	6783323	9600707	1.51%	0.488%
16	12.416	1663	1671	1676	rVB	16579313	26067041	4.10%	1.326%
17	12.627	1696	1707	1712	rBV	9177021	16045985	2.52%	0.816%
18	12.774	1722	1732	1746	rBV2	4215165	16042816	2.52%	0.816%
19	13.104	1777	1788	1793	rBV	47628473	86317195	13.56%	4.390%
20	13.280	1804	1818	1824	rBV	11836622	20078348	3.15%	1.021%
21	13.410	1829	1840	1847	rVB2	5096047	12047611	1.89%	0.613%
22	13.598	1862	1872	1884	rBV	23626519	46346647	7.28%	2.357%
23	13.898	1910	1923	1928	rBV4	2764916	6451561	1.01%	0.328%
24	14.592	2024	2041	2046	rBV	53292698	160114091	25.16%	8.144%
25	14.645	2046	2050	2058	rVB	4406903	7465641	1.17%	0.380%
26	14.798	2072	2076	2080	rVV	5689579	9132006	1.43%	0.464%
27	14.851	2080	2085	2093	rVB2	4813138	8077815	1.27%	0.411%
28	14.974	2100	2106	2110	rBV2	5579471	8856164	1.39%	0.450%
29	15.027	2110	2115	2118	rVV3	4325501	8321344	1.31%	0.423%
30	15.092	2118	2126	2133	rVV2	9834225	29320180	4.61%	1.491%
31	15.157	2133	2137	2141	rVV2	5386474	10067957	1.58%	0.512%
32	15.204	2141	2145	2150	rVB	9388164	13344938	2.10%	0.679%
33	15.321	2155	2165	2182	rVB3	8076303	32631147	5.13%	1.660%
34	15.510	2189	2197	2200	rBV3	3752691	9922328	1.56%	0.505%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	15.562	2200	2206	2213	rVB4	5021311	14236336	2.24%	0.724%
36	15.786	2223	2244	2246	rBV4	136816928	636410036	100.00%	32.370%
37	17.015	2441	2453	2459	rBV	69410275	147900628	23.24%	7.523%
38	17.251	2487	2493	2497	rVV	15936552	20281621	3.19%	1.032%
39	17.386	2511	2516	2521	rVV	7311622	13674513	2.15%	0.696%
40	17.445	2521	2526	2533	rVV3	5813576	11196608	1.76%	0.569%
41	17.515	2533	2538	2548	rVB	6398124	9019016	1.42%	0.459%
42	17.751	2572	2578	2584	rVB	10124672	14762953	2.32%	0.751%

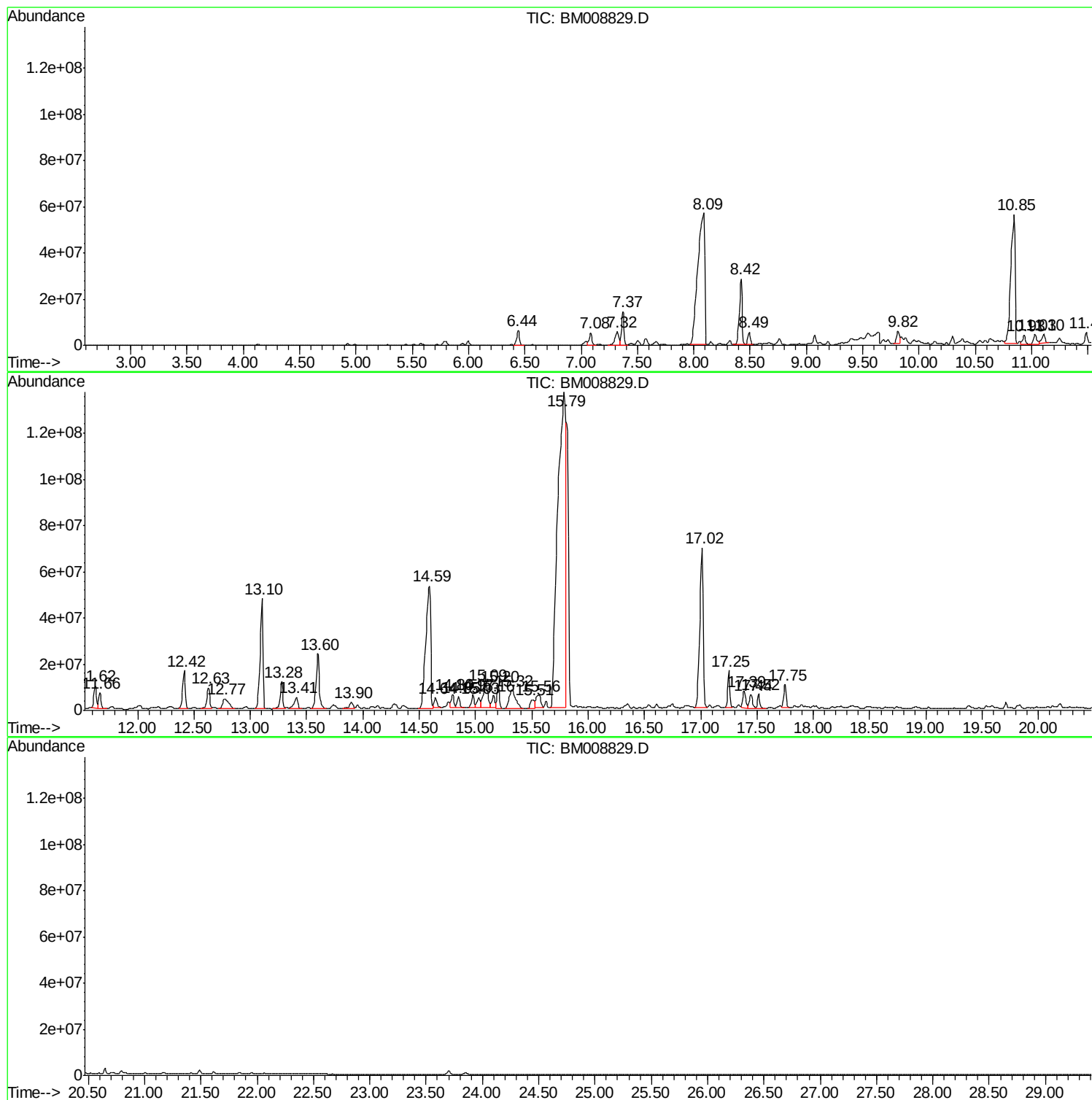
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BNA_M
ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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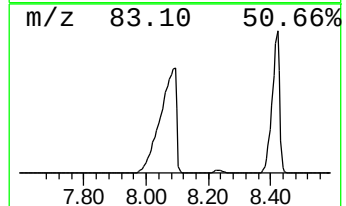
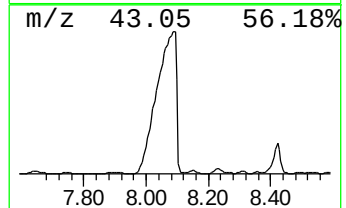
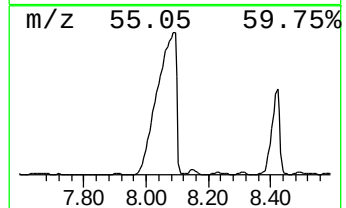
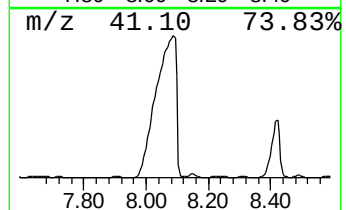
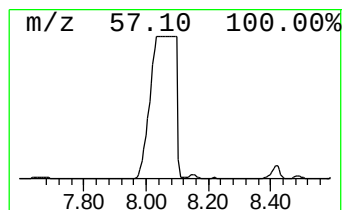
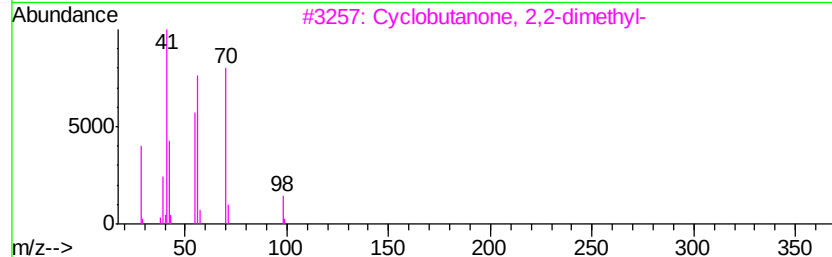
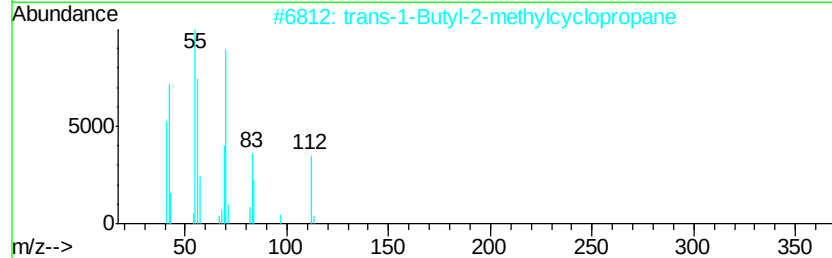
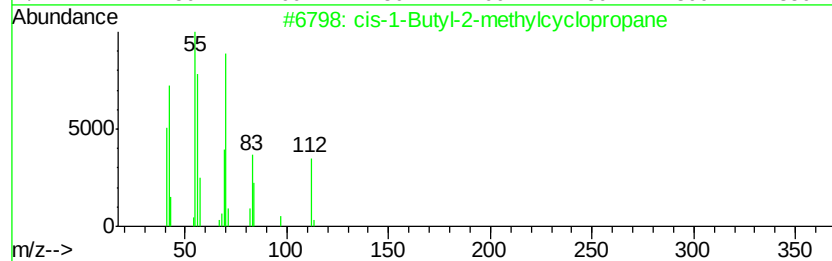
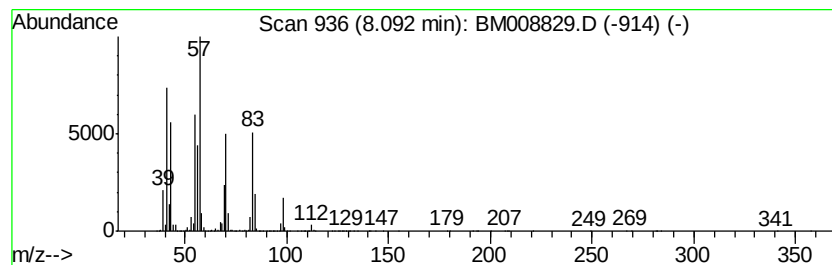
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic8.09 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.09	20.00 ng/ul	244658000	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-1-Butyl-2-methylcyclopropane	112	C8H16	038851-69-3	58
2		trans-1-Butyl-2-methylcyclopropane	112	C8H16	038851-70-6	52
3		Cyclobutanone, 2,2-dimethyl-	98	C6H10O	001192-14-9	43
4		Cyclopropane, 1,1,2,3-tetramethyl-	98	C7H14	074752-93-5	30
5		1-Hexene, 3,5,5-trimethyl-	126	C9H18	004316-65-8	27



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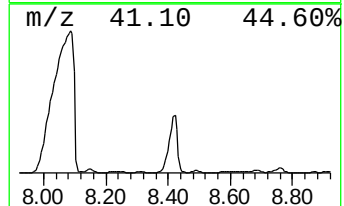
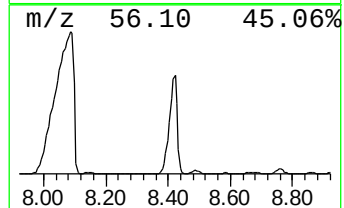
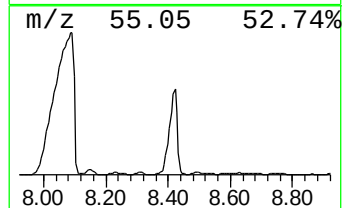
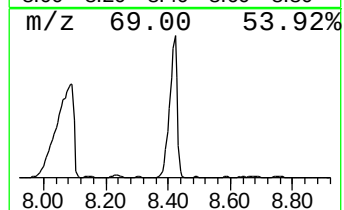
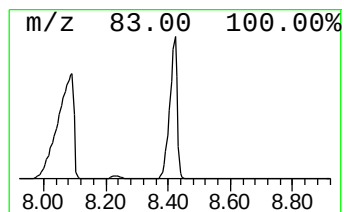
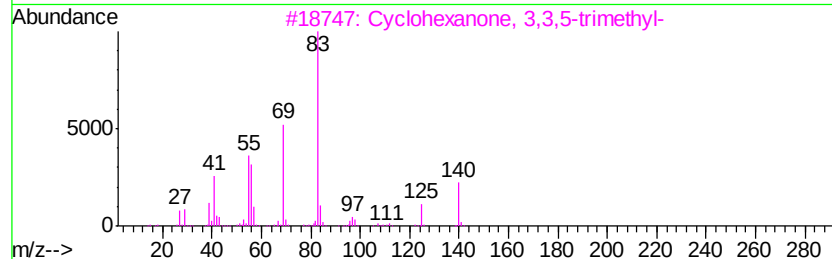
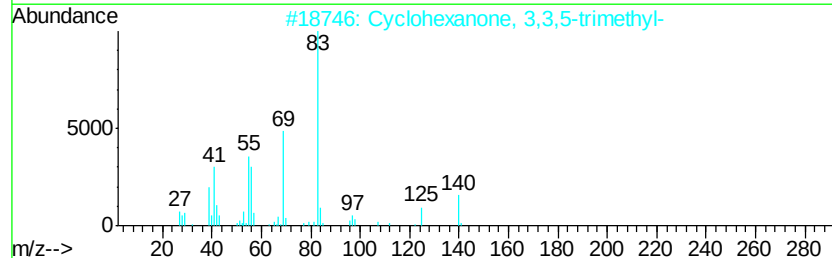
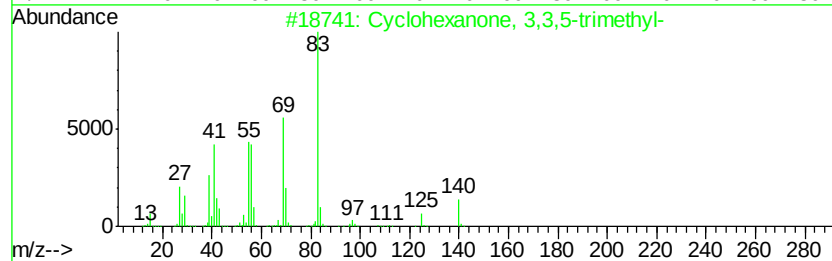
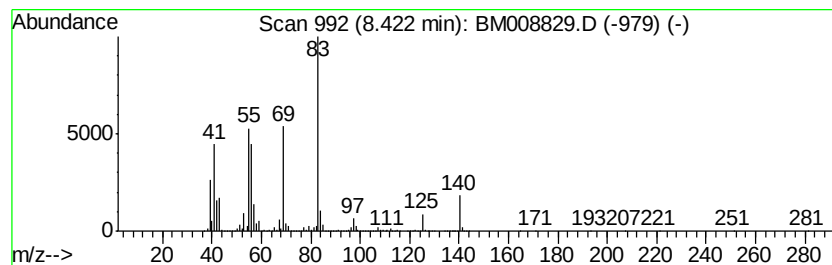
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Cyclohexanone, 3,3,5-trimet... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	3.99 ng/ul	48805300	1,4-Dichlorobenzene-d4	7.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
2		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
3		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	81
4		1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	53
5		Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	53



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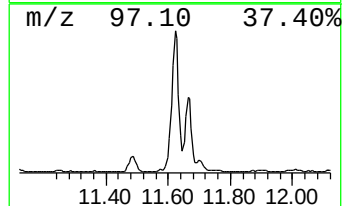
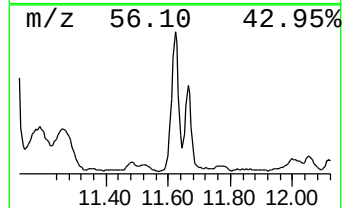
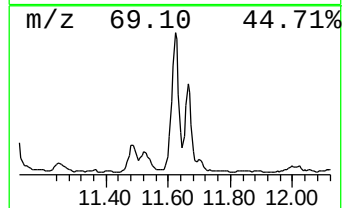
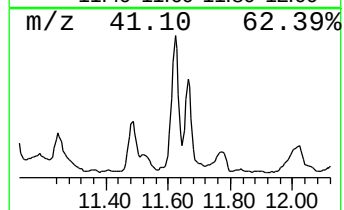
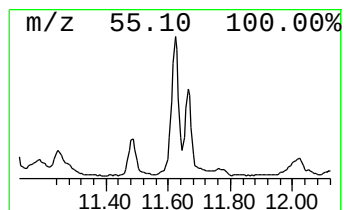
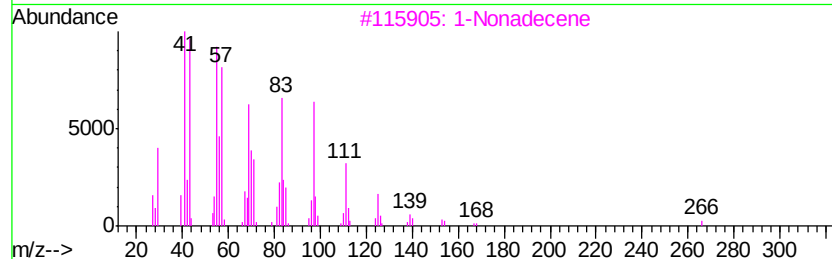
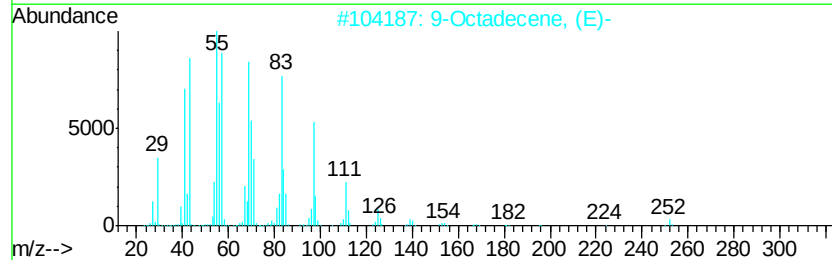
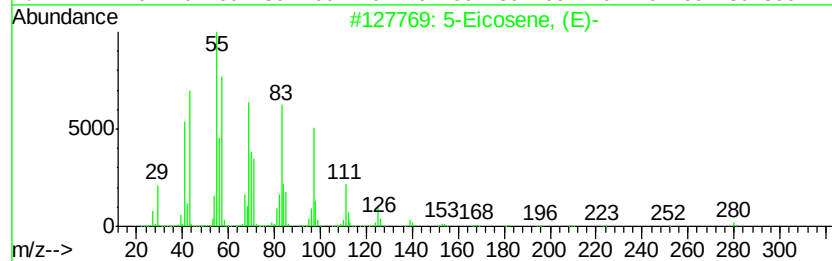
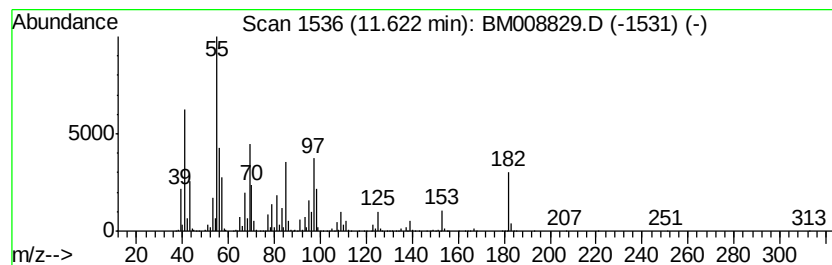
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic11.62 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	2.04 ng/ul	14901800	Napthalene-d8	10.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	37
2		9-Octadecene, (E)-	252	C18H36	007206-25-9	32
3		1-Nonadecene	266	C19H38	018435-45-5	25
4		4-Nonene, 5-butyl-	182	C13H26	007367-38-6	25
5		3-Heptene, (E)-	98	C7H14	014686-14-7	22



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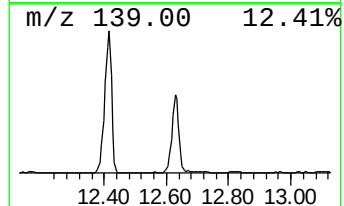
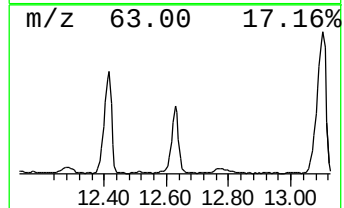
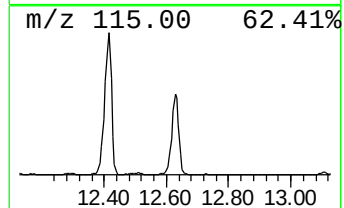
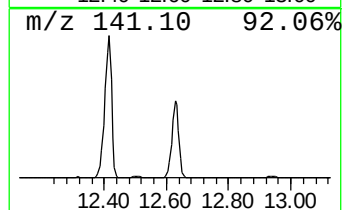
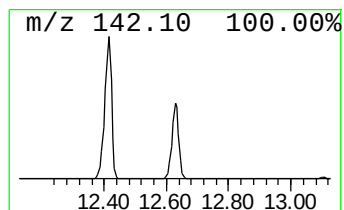
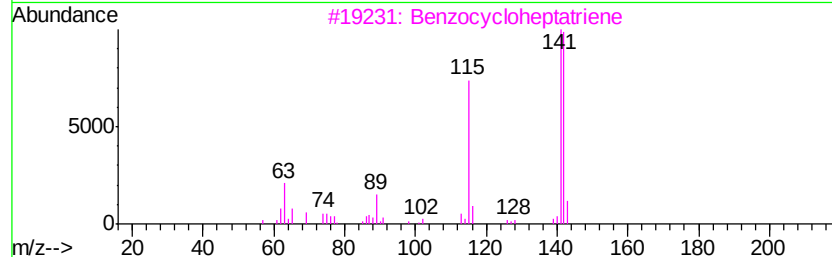
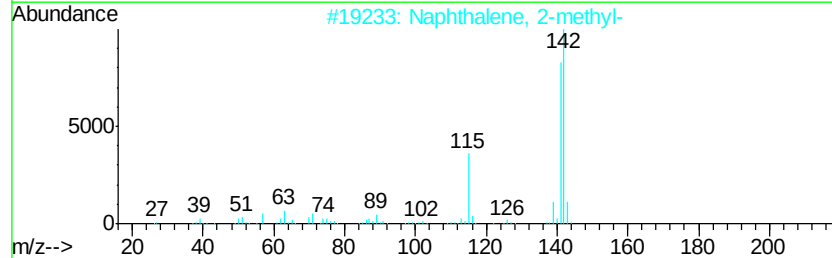
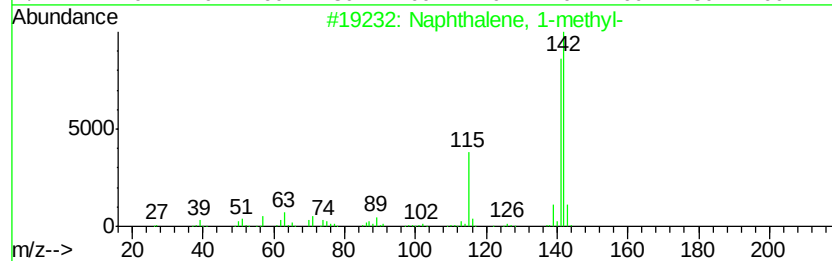
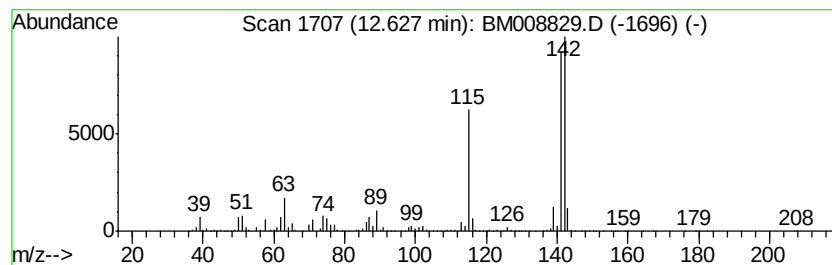
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Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	2.20 ng/ul	16046000	Naphthalene-d8	10.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
3		Benzocycloheptatriene	142	C11H10	000264-09-5	96
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94



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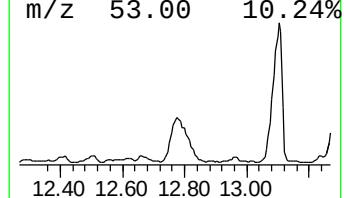
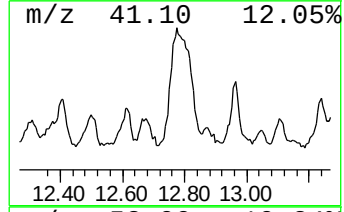
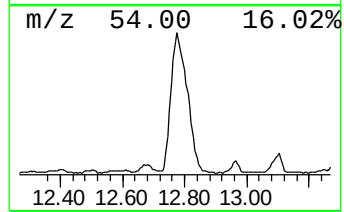
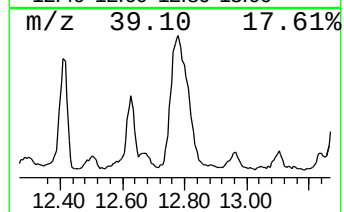
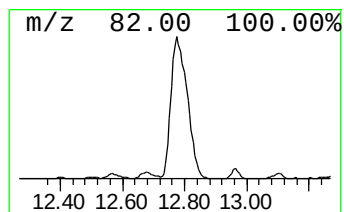
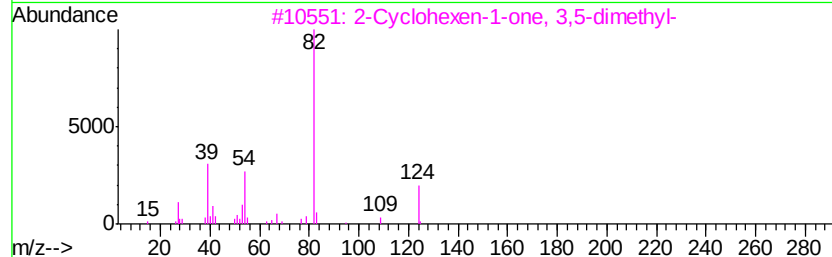
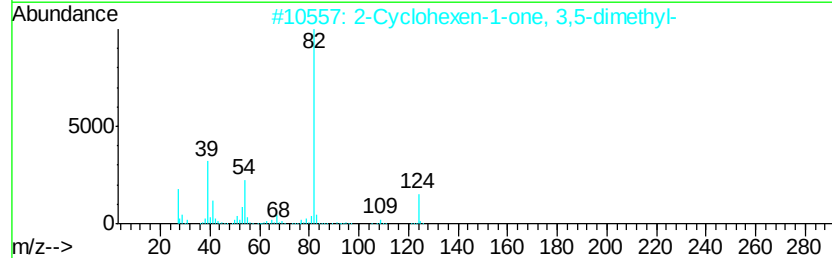
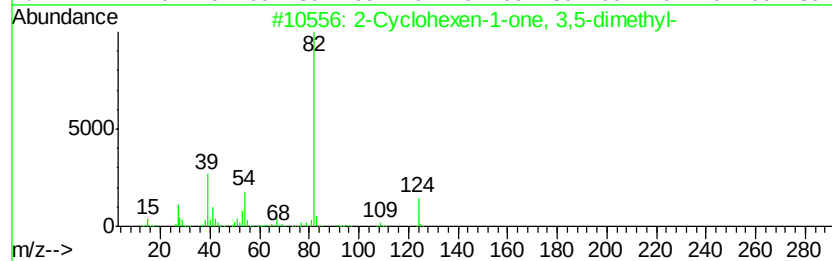
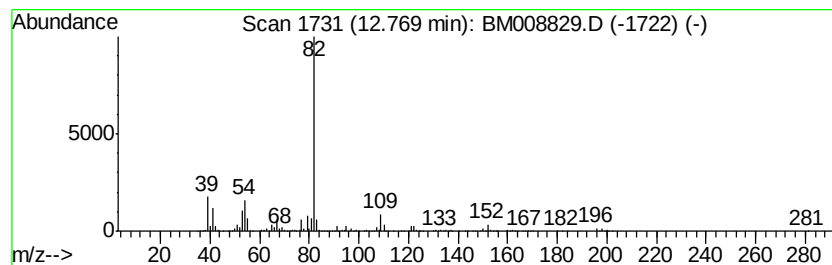
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Cyclohexen-1-one, 3,5-dim... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	2.00 ng/ul	16042800	Acenaphthene-d10	14.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Cyclohexen-1-one, 3,5-dimethyl-	124	C8H12O	001123-09-7	64
2		2-Cyclohexen-1-one, 3,5-dimethyl-	124	C8H12O	001123-09-7	64
3		2-Cyclohexen-1-one, 3,5-dimethyl-	124	C8H12O	001123-09-7	59
4		1H-Pyrazole, 1-methyl-	82	C4H6N2	000930-36-9	50
5		Isophorone	138	C9H14O	000078-59-1	50



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
Data File : BM008829.D
Acq On : 24 Jan 2017 07:44
Operator : UM/SJ
Sample : I1323-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

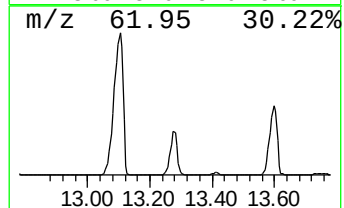
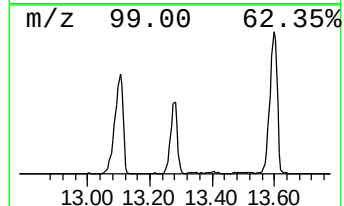
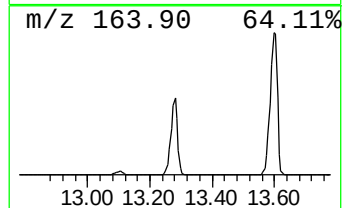
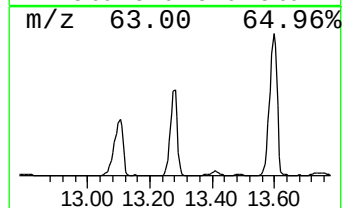
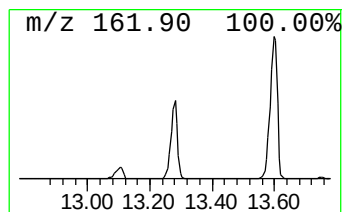
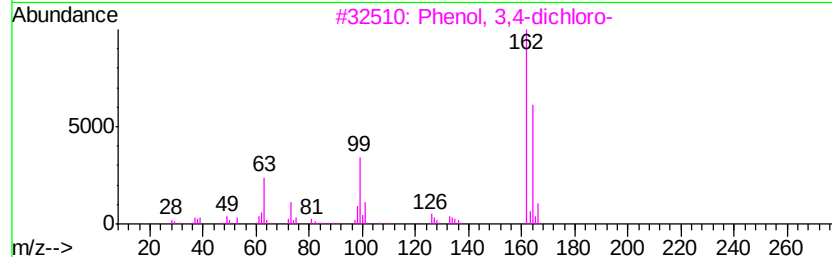
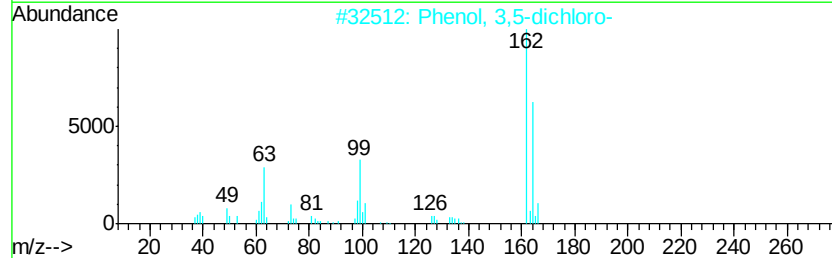
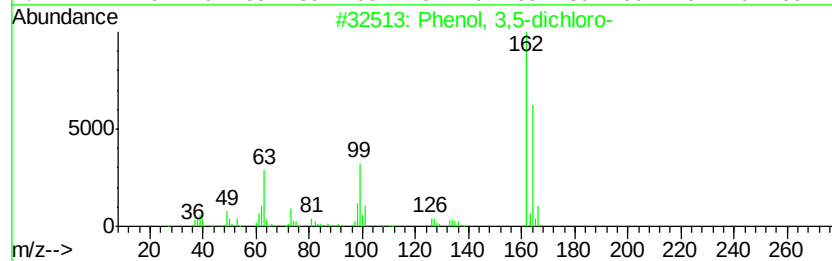
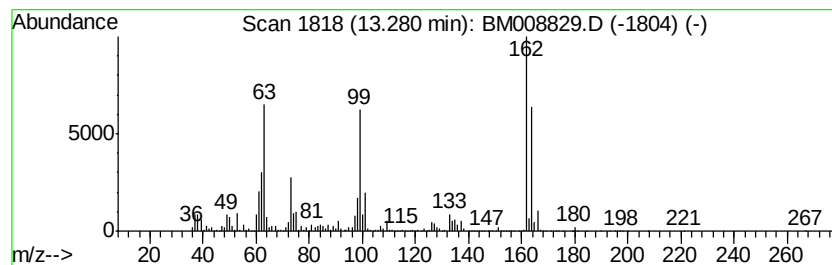
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 3,5-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.28	2.51 ng/ul	20078300	Acenaphthene-d10	14.58		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol,	3,5-dichloro-	162	C6H4Cl2O	000591-35-5	93
2	Phenol,	3,5-dichloro-	162	C6H4Cl2O	000591-35-5	93
3	Phenol,	3,4-dichloro-	162	C6H4Cl2O	000095-77-2	91
4	Phenol,	3,4-dichloro-	162	C6H4Cl2O	000095-77-2	91
5	Phenol,	3,5-dichloro-	162	C6H4Cl2O	000591-35-5	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
Data File : BM008829.D
Acq On : 24 Jan 2017 07:44
Operator : UM/SJ
Sample : I1323-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

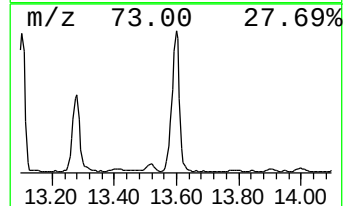
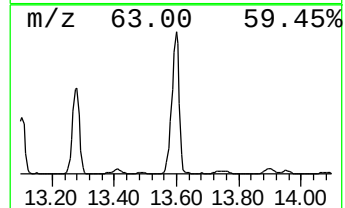
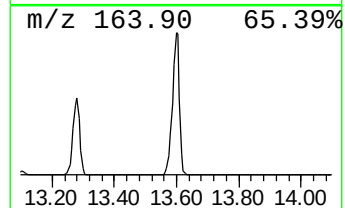
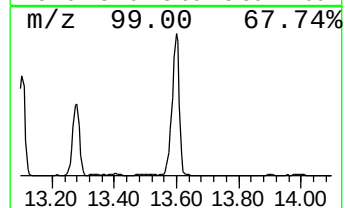
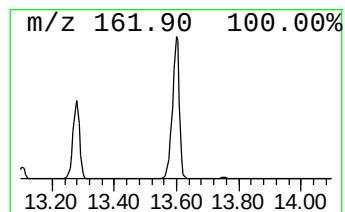
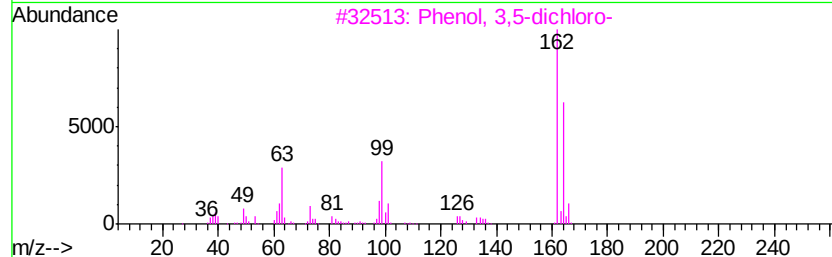
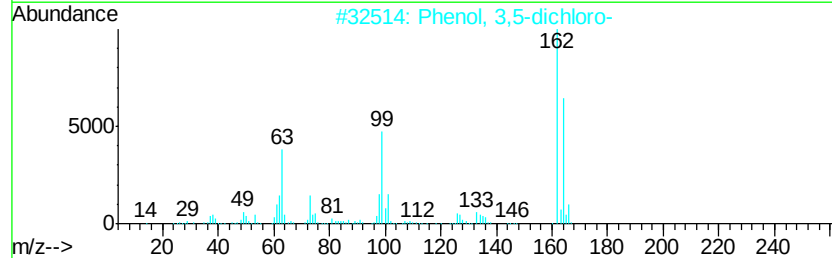
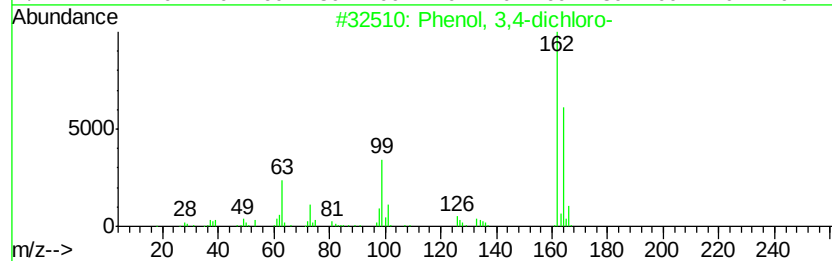
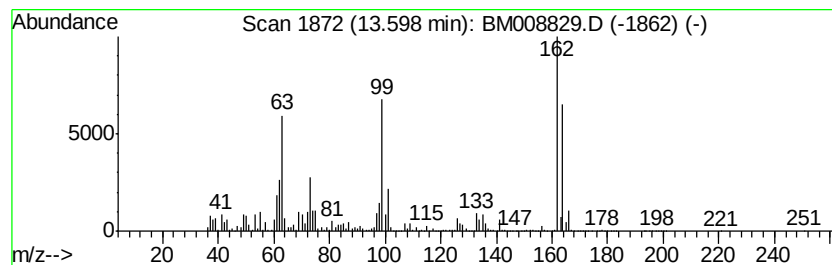
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, 3,4-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	5.79 ng/ul	46346600	Acenaphthene-d10	14.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2 97
2	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5 94
3	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5 91
4	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2 91
5	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5 91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
Data File : BM008829.D
Acq On : 24 Jan 2017 07:44
Operator : UM/SJ
Sample : I1323-22
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ALS Vial : 29 Sample Multiplier: 1

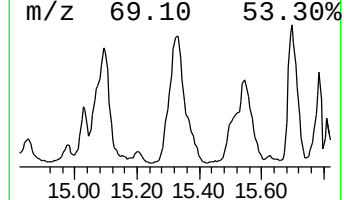
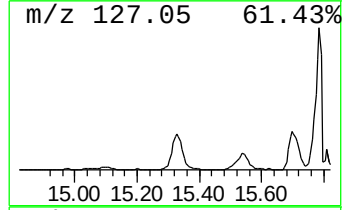
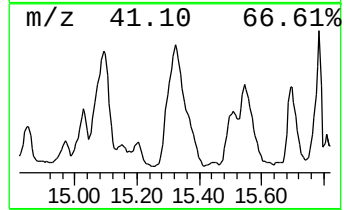
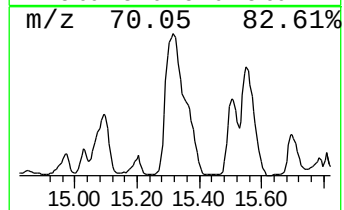
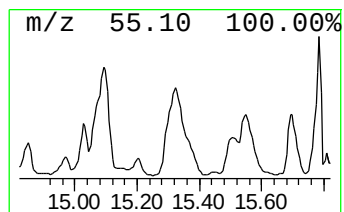
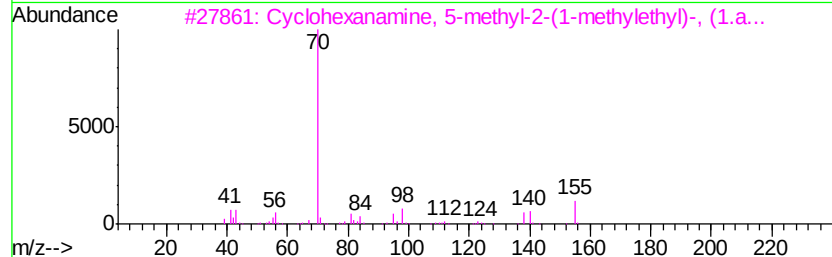
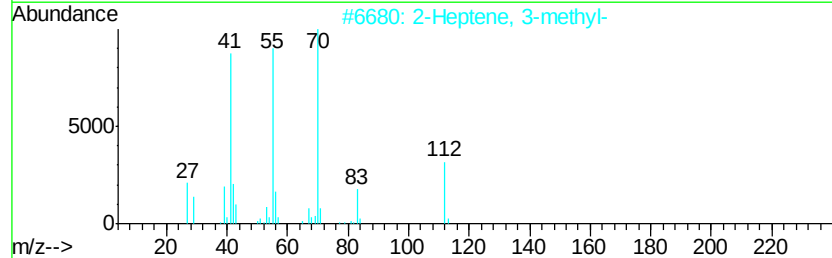
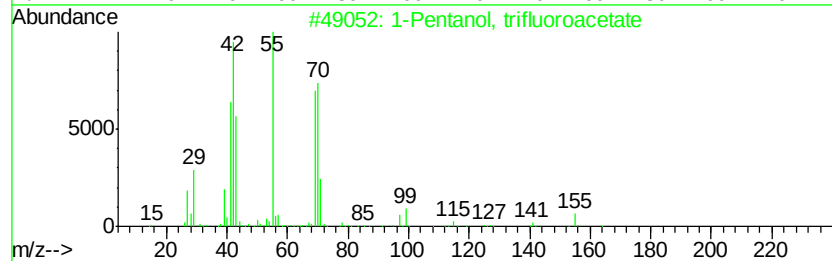
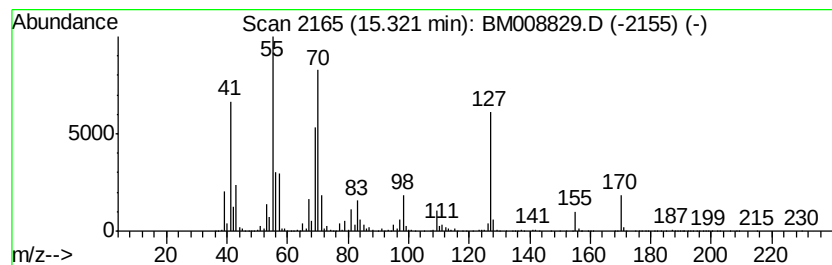
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.32	4.08 ng/ul	32631100	Acenaphthene-d10	14.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentanol, trifluoroacetate	184	C7H11F3O2	000327-70-8	43
2	2-Heptene, 3-methyl-	112	C8H16	003404-75-9	38
3	Cyclohexanamine, 5-methyl-2-(1-m...	155	C10H21N	023399-21-5	38
4	Cycloheptane	98	C7H14	000291-64-5	38
5	Cyclopentane, 1,1,3,3-tetramethyl-	126	C9H18	050876-33-0	38



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
Data File : BM008829.D
Acq On : 24 Jan 2017 07:44
Operator : UM/SJ
Sample : I1323-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

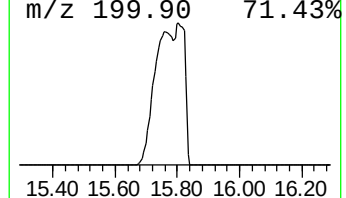
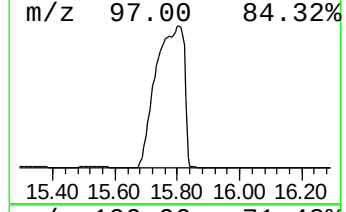
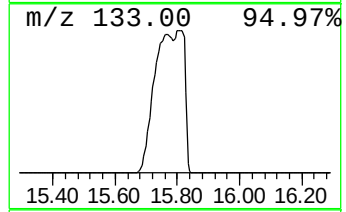
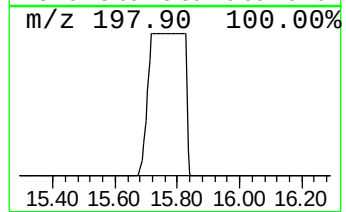
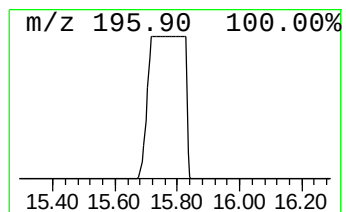
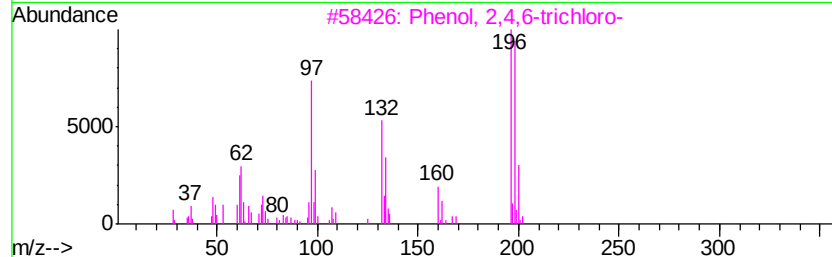
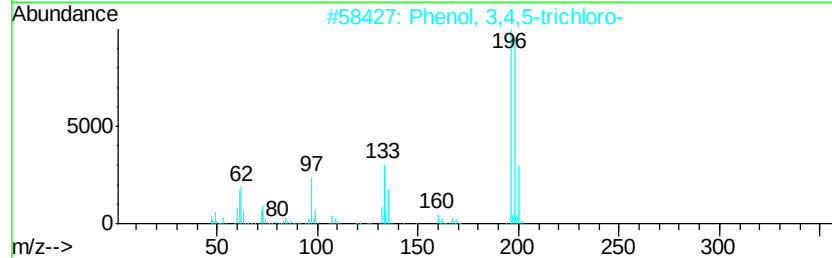
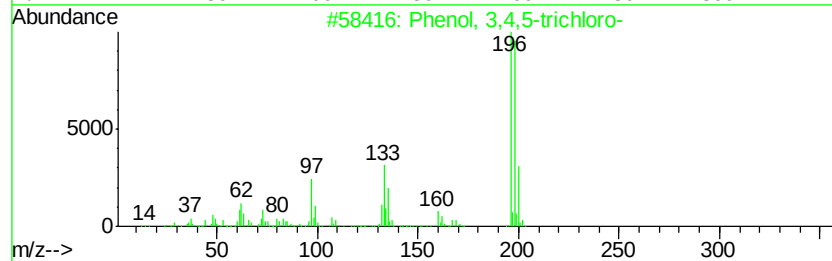
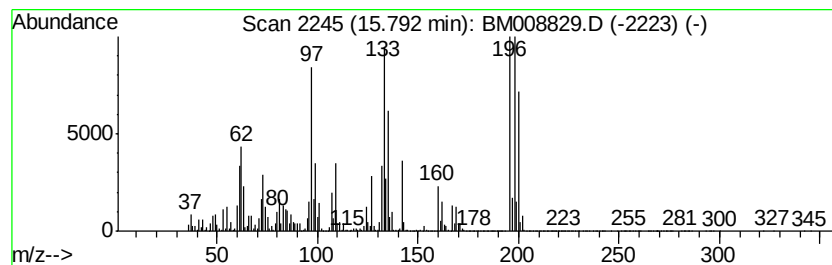
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phenol, 3,4,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	79.49 ng/ul	636410000	Acenaphthene-d10	14.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, 3,4,5-trichloro-	196 C6H3Cl3O	000609-19-8	91
2	Phenol, 3,4,5-trichloro-	196 C6H3Cl3O	000609-19-8	90
3	Phenol, 2,4,6-trichloro-	196 C6H3Cl3O	000088-06-2	83
4	Phenol, 2,4,6-trichloro-	196 C6H3Cl3O	000088-06-2	64
5	Phenol, 2,3,5-trichloro-	196 C6H3Cl3O	000933-78-8	64



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
Data File : BM008829.D
Acq On : 24 Jan 2017 07:44
Operator : UM/SJ
Sample : I1323-22
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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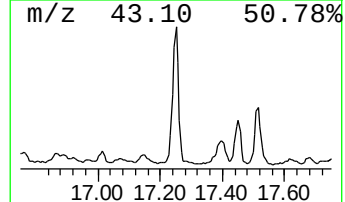
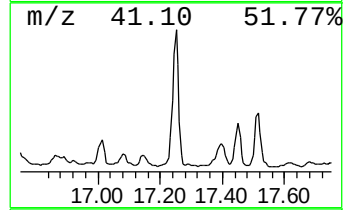
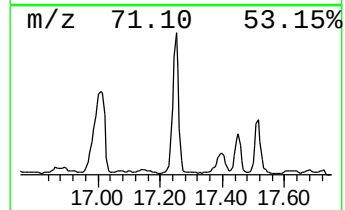
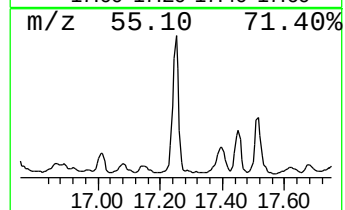
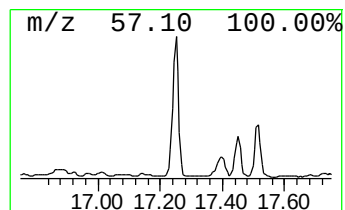
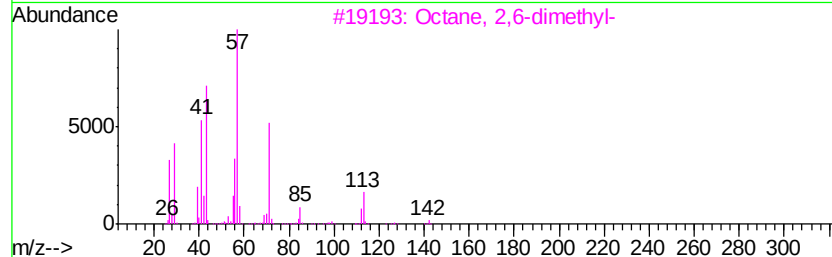
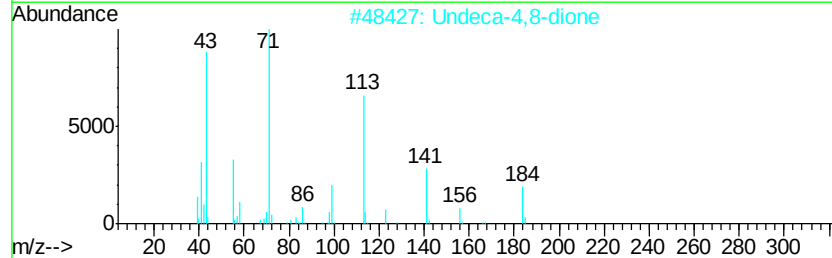
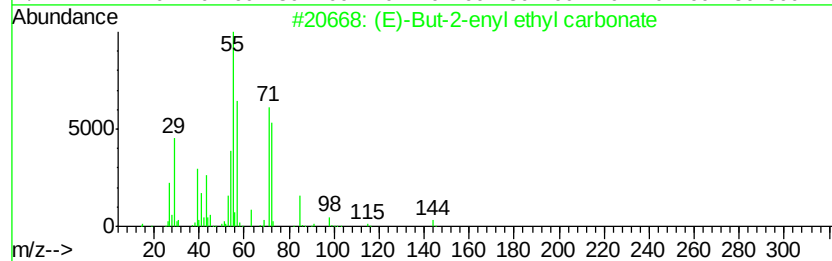
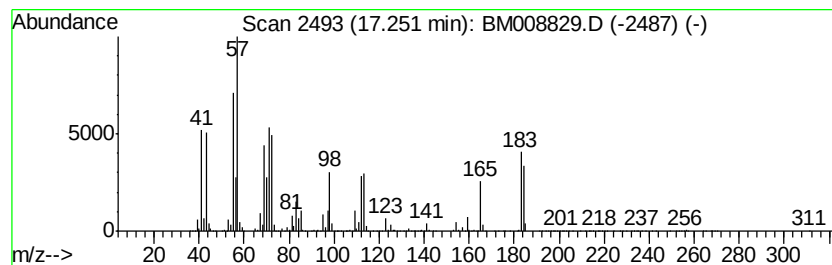
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.25	29.66 ng/ul	20281600	Phenanthrene-d10	17.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-But-2-enyl ethyl carbonate	144	C7H12O3	1000373-77-0	22
2		Undeca-4,8-dione	184	C11H20O2	013505-35-6	22
3		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	14
4		Heptane, 5-ethyl-2,2,3-trimethyl-	170	C12H26	062199-06-8	14
5		Sulfurous acid, 2-ethylhexyl oct...	446	C26H54O3S	1000309-20-1	14



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012317\
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Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DA9R7

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012317.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Cyc...	8.09	20.0	ng/ul	244658000	1	7.94	244658000	20.0
Cyclohexanone, 3,...	8.42	4.0	ng/ul	48805300	1	7.94	244658000	20.0
(DEL) Alkane: Cyc...	11.62	2.0	ng/ul	14901800	2	10.76	146016000	20.0
Naphthalene, 1-me...	12.63	2.2	ng/ul	16046000	2	10.76	146016000	20.0
2-Cyclohexen-1-on...	12.77	2.0	ng/ul	16042800	3	14.58	160114000	20.0
Phenol, 3,5-dichl...	13.28	2.5	ng/ul	20078300	3	14.58	160114000	20.0
Phenol, 3,4-dichl...	13.60	5.8	ng/ul	46346600	3	14.58	160114000	20.0
unknown-01	15.32	4.1	ng/ul	32631100	3	14.58	160114000	20.0
Phenol, 3,4,5-tri...	15.79	79.5	ng/ul	636410000	3	14.58	160114000	20.0
unknown-02	17.25	29.7	ng/ul	20281600	4	17.34	13674500	20.0