

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
 Data File : BN004316.D
 Acq On : 03 Jan 2019 13:58
 Operator : JU/SJ
 Sample : J6593-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BE803

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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.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	7.146	701	707	711	rBV	109497	167200	2.35%	0.686%
2	7.340	734	740	745	rBV	124258	193844	2.73%	0.795%
3	7.810	814	820	826	rVB	151221	234748	3.30%	0.963%
4	8.169	875	881	887	rBV	58432	96187	1.35%	0.395%
5	8.516	936	940	952	rVB	94913	152921	2.15%	0.627%
6	8.981	1013	1019	1025	rBV	161400	256032	3.60%	1.050%
7	9.269	1056	1068	1076	rBV	157461	298975	4.21%	1.227%
8	9.704	1087	1142	1144	rBV3	816043	7108377	100.00%	29.165%
9	10.163	1211	1220	1224	rBV	47955	126691	1.78%	0.520%
10	10.234	1224	1232	1236	rVV2	255813	511187	7.19%	2.097%
11	10.281	1236	1240	1252	rVB2	196039	376013	5.29%	1.543%
12	10.546	1277	1285	1289	rBV2	64585	128289	1.80%	0.526%
13	10.604	1289	1295	1302	rVB	228384	392929	5.53%	1.612%
14	10.787	1316	1326	1331	rBV5	33083	87500	1.23%	0.359%
15	10.899	1339	1345	1348	rBV	55326	111471	1.57%	0.457%
16	10.957	1348	1355	1358	rBV	103652	193144	2.72%	0.792%
17	11.104	1372	1380	1389	rVB2	177512	424088	5.97%	1.740%
18	11.193	1389	1395	1398	rBV4	47697	91521	1.29%	0.376%
19	11.240	1398	1403	1406	rBV4	58112	109327	1.54%	0.449%
20	11.404	1425	1431	1435	rBV	85289	206371	2.90%	0.847%
21	11.551	1448	1456	1461	rVB3	201375	521745	7.34%	2.141%
22	11.616	1462	1467	1470	rBV	56180	89455	1.26%	0.367%
23	11.669	1470	1476	1479	rBV2	122510	246543	3.47%	1.012%
24	11.775	1488	1494	1498	rBV2	146167	276130	3.88%	1.133%
25	11.846	1498	1506	1510	rVV	181361	369976	5.20%	1.518%
26	11.887	1510	1513	1519	rVB	185397	268419	3.78%	1.101%
27	12.028	1529	1537	1538	rBV3	60559	117081	1.65%	0.480%
28	12.063	1538	1543	1548	rVV	255422	393252	5.53%	1.613%
29	12.116	1548	1552	1561	rVB	296222	542756	7.64%	2.227%
30	12.298	1577	1583	1590	rBV3	202023	405982	5.71%	1.666%
31	12.381	1593	1597	1601	rBV4	45375	72265	1.02%	0.296%
32	12.451	1606	1609	1617	rVB3	79563	161856	2.28%	0.664%
33	12.610	1629	1636	1647	rVB3	173322	411960	5.80%	1.690%
34	12.710	1647	1653	1660	rBV2	116918	268267	3.77%	1.101%

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35	12.928	1681	1690	1694	rBV2	246255	500544	7.04%	2.054%
36	12.969	1694	1697	1701	rVV3	73110	105912	1.49%	0.435%
37	13.016	1701	1705	1712	rVV5	80258	193957	2.73%	0.796%
38	13.140	1724	1726	1732	rVB2	72677	100426	1.41%	0.412%
39	13.216	1735	1739	1751	rBV3	51430	184565	2.60%	0.757%
40	13.334	1755	1759	1761	rBV3	50246	80298	1.13%	0.329%
41	13.387	1762	1768	1773	rVB5	145023	302205	4.25%	1.240%
42	13.628	1805	1809	1814	rBV3	65544	130061	1.83%	0.534%
43	13.716	1820	1824	1827	rBV4	100790	181073	2.55%	0.743%
44	13.863	1844	1849	1853	rBV	299797	474157	6.67%	1.945%
45	13.951	1861	1864	1871	rVB2	119513	220874	3.11%	0.906%
46	14.104	1886	1890	1893	rBV3	160721	281993	3.97%	1.157%
47	14.151	1893	1898	1903	rVB	408027	715002	10.06%	2.934%
48	14.328	1925	1928	1932	rBV4	82676	135115	1.90%	0.554%
49	14.457	1945	1950	1958	rBV3	291466	568396	8.00%	2.332%
50	15.016	2042	2045	2049	rVB3	101090	141120	1.99%	0.579%
51	15.445	2113	2118	2122	rBV	448443	610399	8.59%	2.504%
52	15.575	2135	2140	2147	rVB3	282547	562289	7.91%	2.307%
53	17.198	2412	2416	2425	rVB2	334386	537038	7.56%	2.203%
54	17.298	2428	2433	2440	rVB2	438584	647585	9.11%	2.657%
55	19.598	2819	2824	2838	rVB2	491455	664937	9.35%	2.728%
56	21.386	3123	3128	3133	rBV	382302	490073	6.89%	2.011%
57	23.557	3491	3497	3504	rVB	316237	595551	8.38%	2.443%
58	23.704	3515	3522	3533	rVB2	245369	460748	6.48%	1.890%
59	26.074	3920	3925	3935	rVB2	31442	76109	1.07%	0.312%

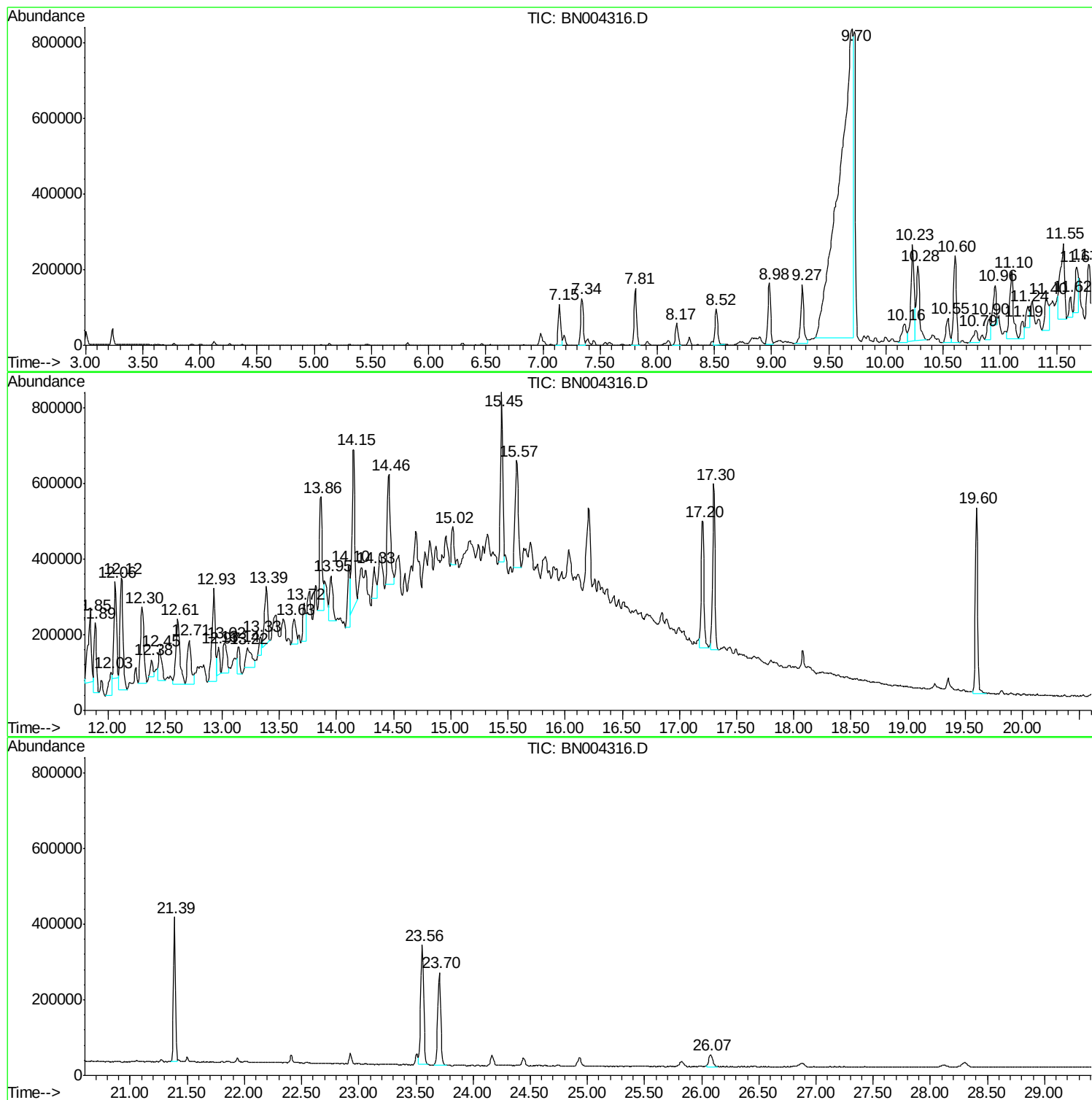
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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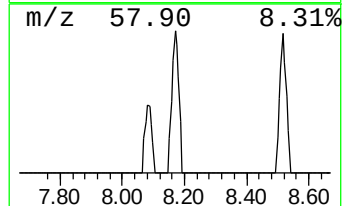
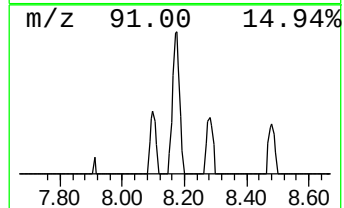
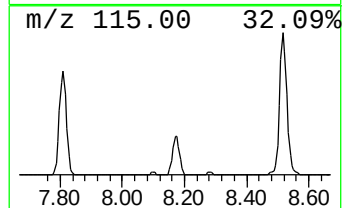
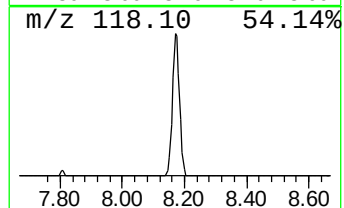
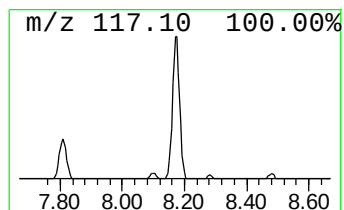
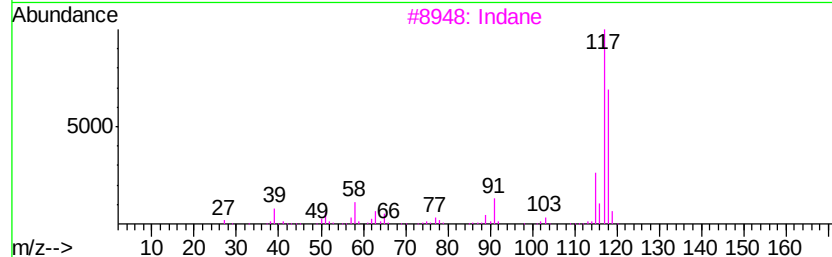
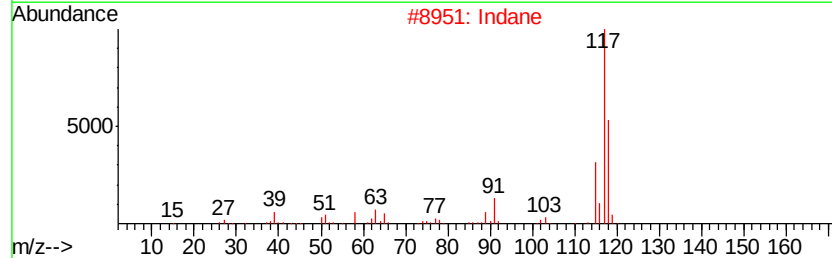
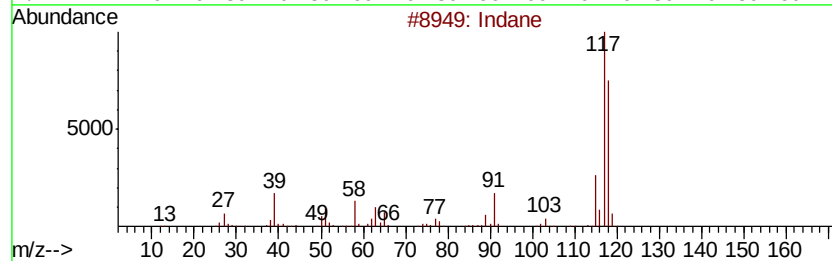
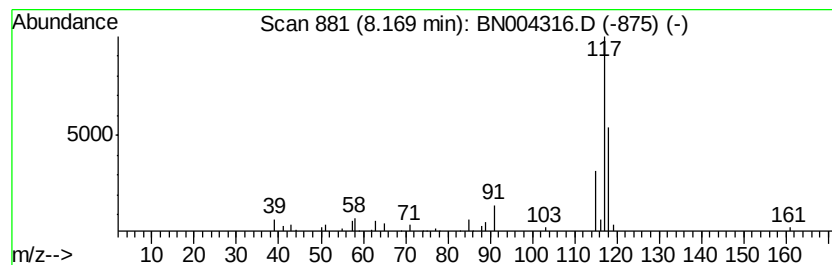
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	8.19 ng/ul	96187	1,4-Dichlorobenzene-d4	7.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	81
2	Indane			118	C9H10	000496-11-7	74
3	Indane			118	C9H10	000496-11-7	72
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-			118	C9H10	1000191-13-7	72
5	Benzene, 2-propenyl-			118	C9H10	000300-57-2	59



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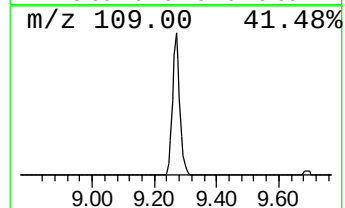
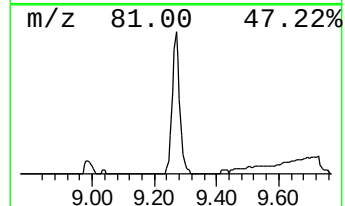
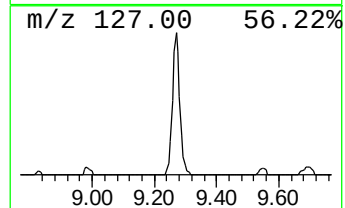
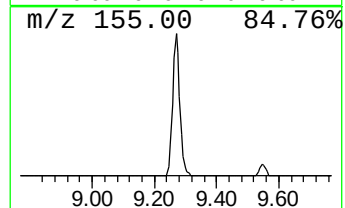
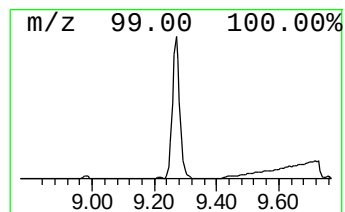
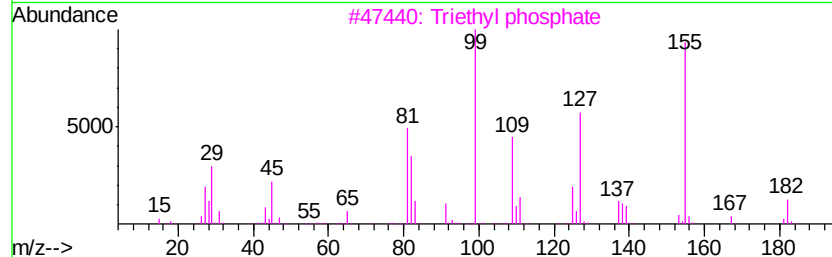
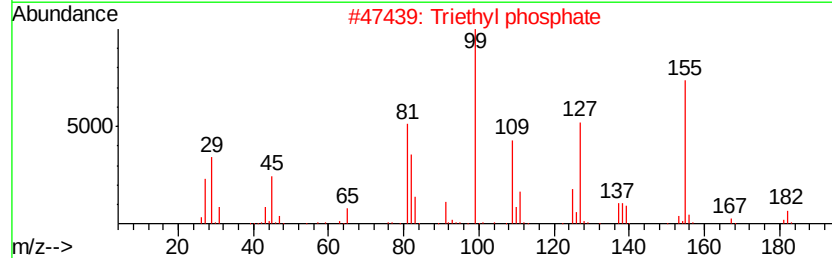
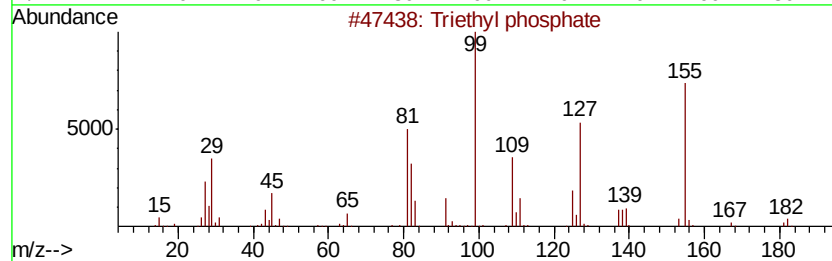
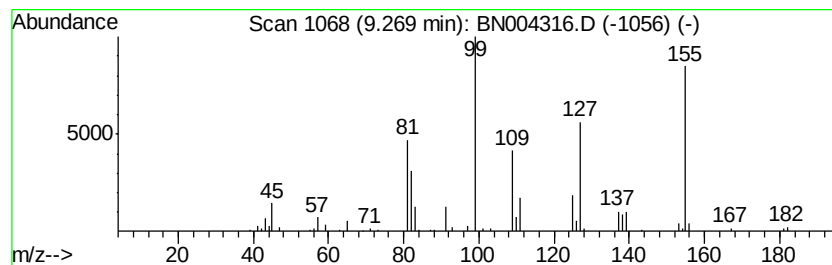
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Triethyl phosphate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.27	15.22 ng/ul	298975	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triethyl phosphate	182	C6H15O4P	000078-40-0	94
2		Triethyl phosphate	182	C6H15O4P	000078-40-0	91
3		Triethyl phosphate	182	C6H15O4P	000078-40-0	90
4		Phosphoric acid, diethyl octyl e...	266	C12H27O4P	1000308-90-4	53
5		Diethyl isopropylidenemalonate	200	C10H16O4	006802-75-1	37



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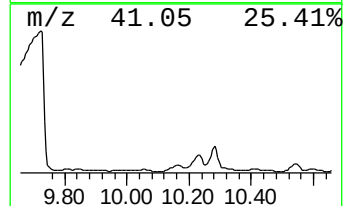
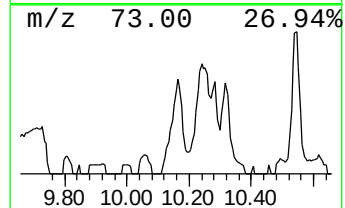
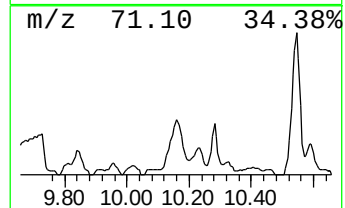
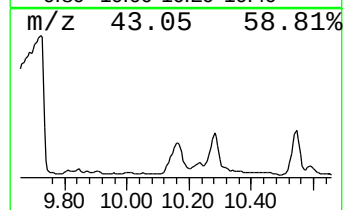
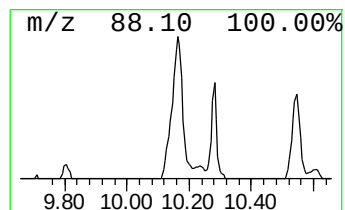
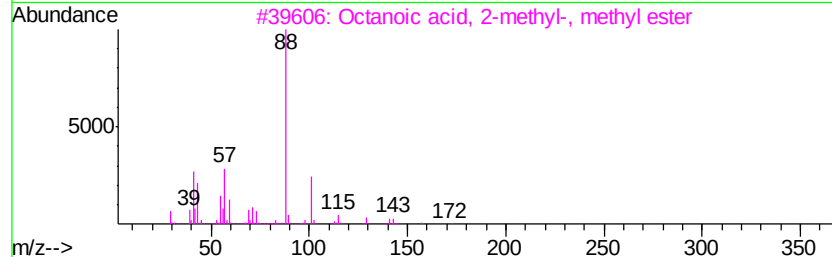
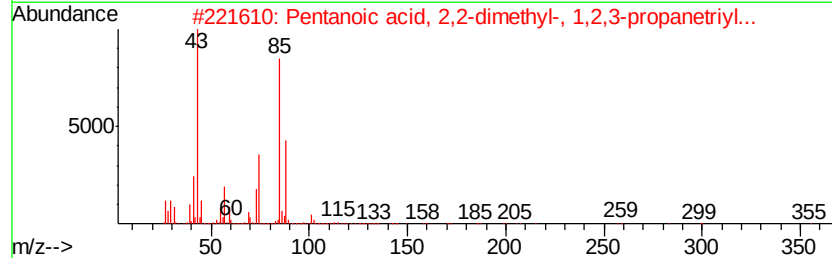
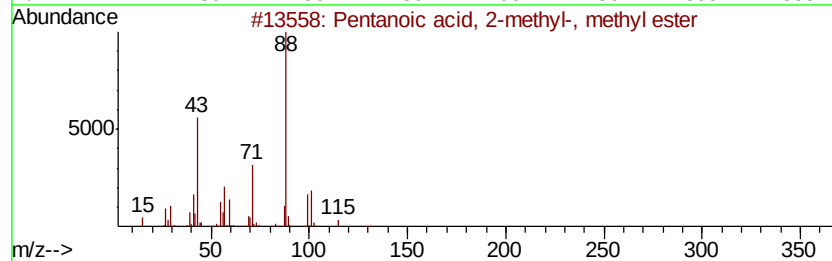
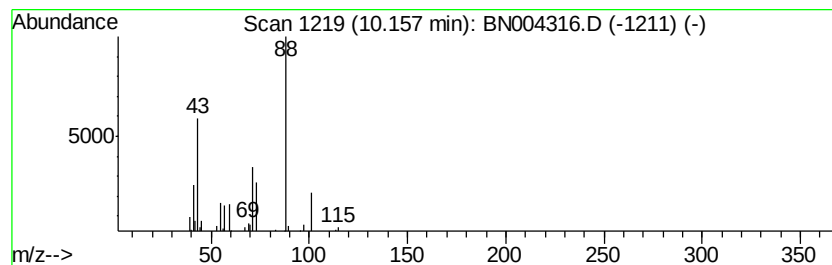
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Pentanoic acid, 2-methyl-, ... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.16	6.45 ng/ul	126691	Napthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentanoic acid, 2-methyl-, methy...	130	C7H14O2	002177-77-7	59
2		Pentanoic acid, 2,2-dimethyl-, 1...	428	C24H44O6	057346-62-0	53
3		Octanoic acid, 2-methyl-, methyl...	172	C10H20O2	002177-86-8	42
4		Pentanoic acid, 2,4-dimethyl-, m...	144	C8H16O2	071672-33-8	42
5		Pentanoic acid, 2-methyl-, methy...	130	C7H14O2	002177-77-7	42



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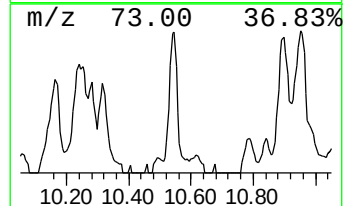
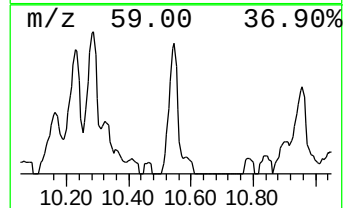
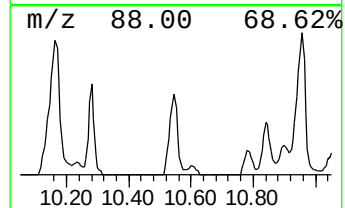
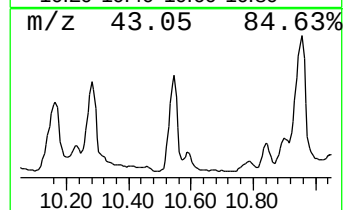
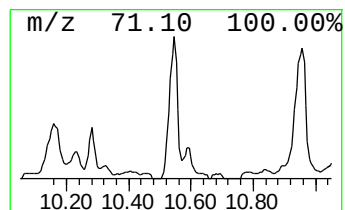
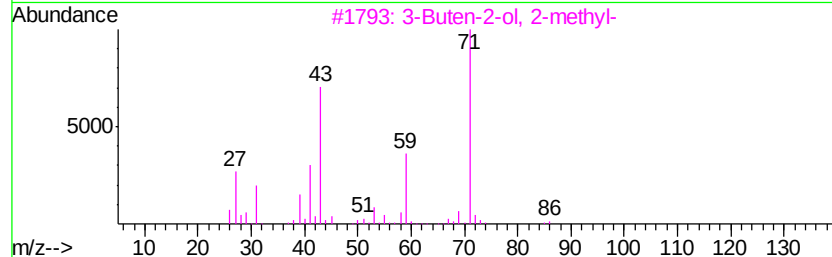
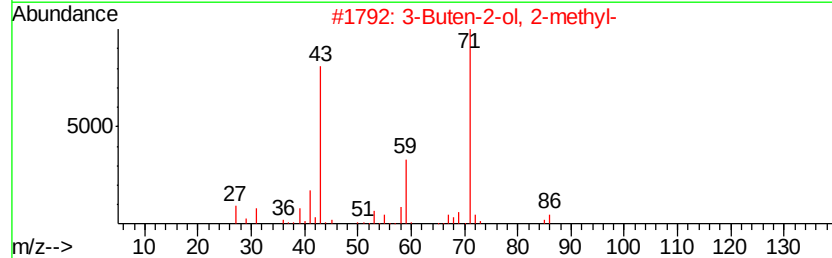
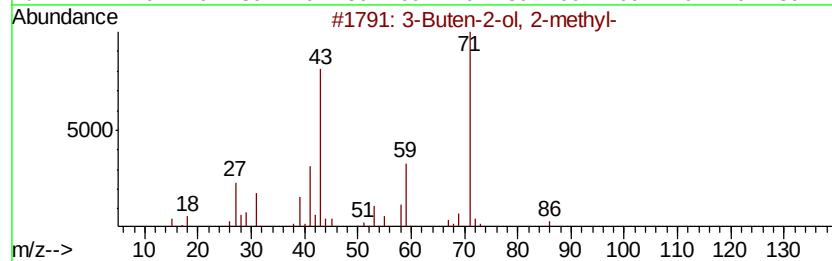
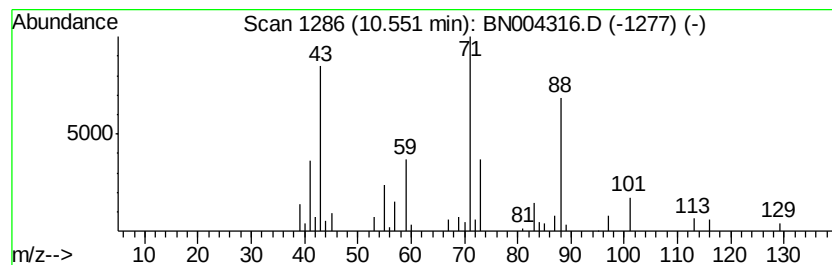
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Peak Number 4 unknown-01 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.55	6.53 ng/ul	128289	Napthalene-d8	10.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	43
2	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	43
3	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	43
4	Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	38
5	2-Ethoxytetrahydrofuran	116	C6H12O2	013436-46-9	35



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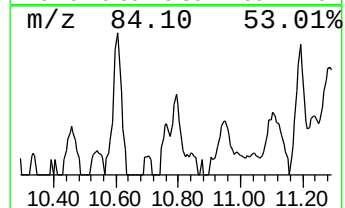
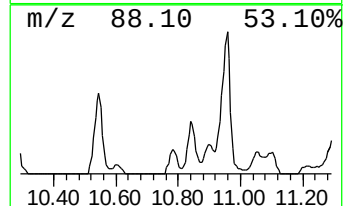
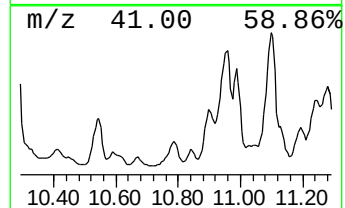
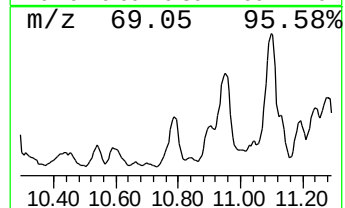
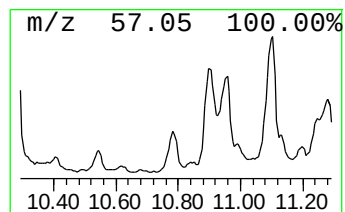
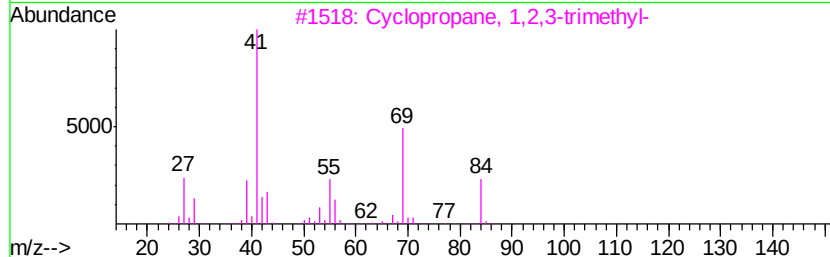
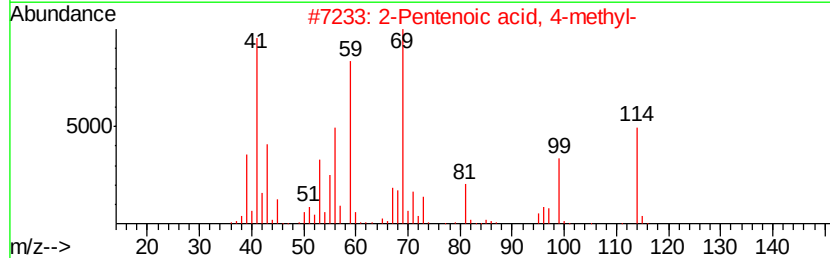
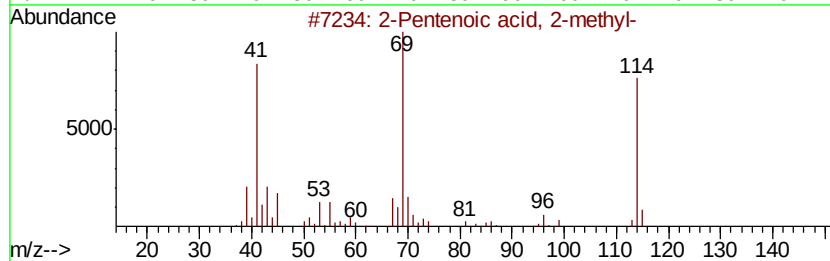
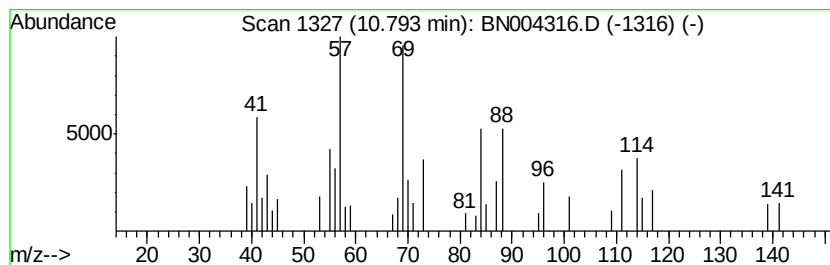
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	4.45 ng/ul	87500	Napthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentenoic acid, 2-methyl-	114	C6H10O2	003142-72-1	27
2		2-Pentenoic acid, 4-methyl-	114	C6H10O2	010321-71-8	27
3		Cyclopropane, 1,2,3-trimethyl-	84	C6H12	042984-19-0	25
4		2-Pentene, 3-methyl-	84	C6H12	000922-61-2	25
5		3-Hexanol, 2,5-dimethyl-	130	C8H18O	019550-07-3	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

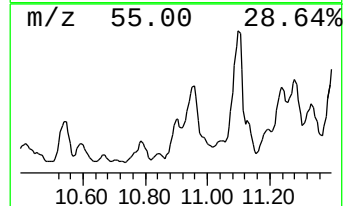
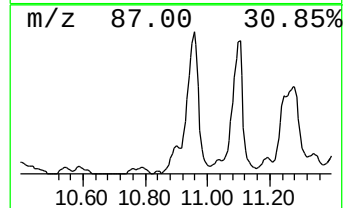
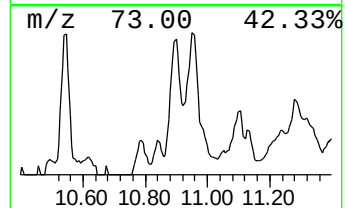
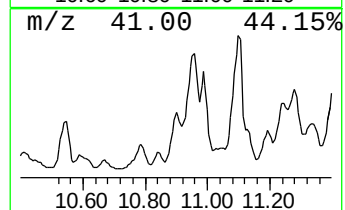
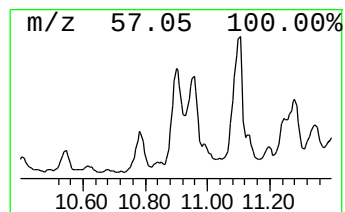
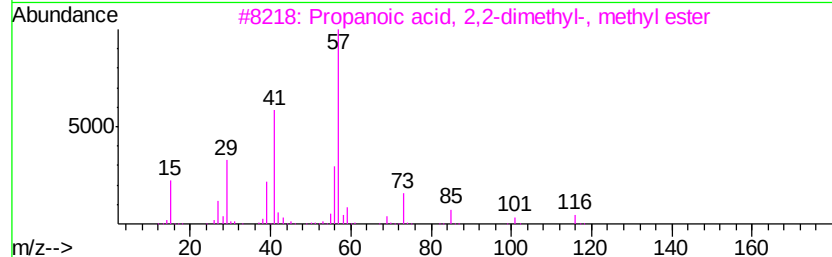
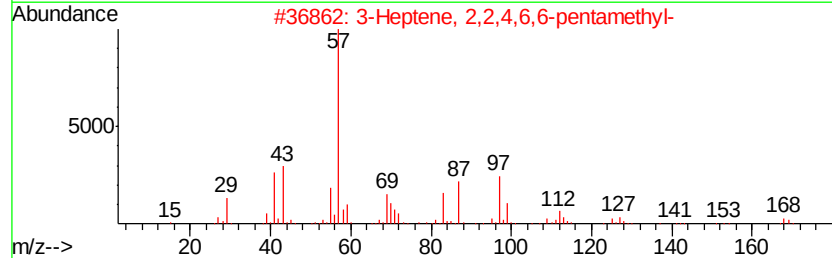
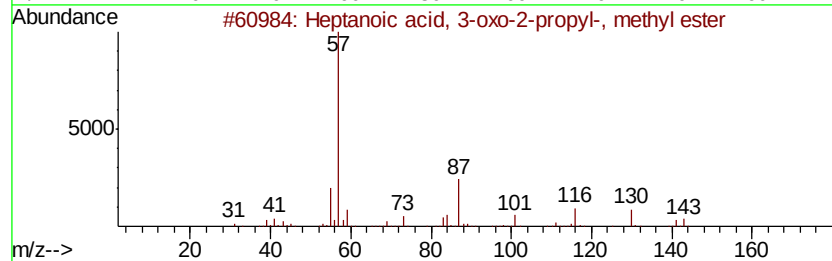
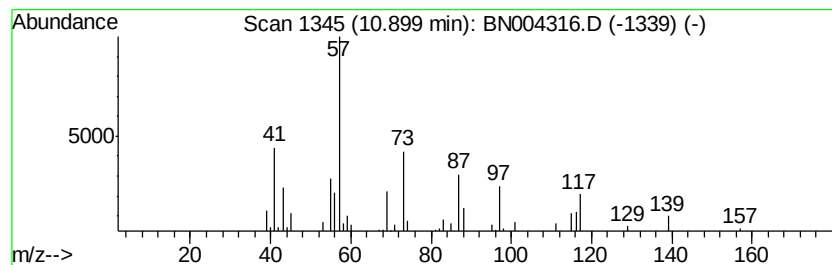
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-03 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.90	5.67 ng/ul	111471	Naphthalene-d8	10.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid, 3-oxo-2-propyl-,...	200	C11H20O3	125055-03-0	35
2	3-Heptene, 2,2,4,6,6-pentamethyl-	168	C12H24	000123-48-8	25
3	Propanoic acid, 2,2-dimethyl-, m...	116	C6H12O2	000598-98-1	25
4	2-Oxepanol, 5-(1,1-dimethylethyl)-	172	C10H20O2	111711-01-4	25
5	Butoxyacetic acid	132	C6H12O3	002516-93-0	17



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
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Sample : J6593-01
Misc :
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Instrument :
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ClientSampleId :
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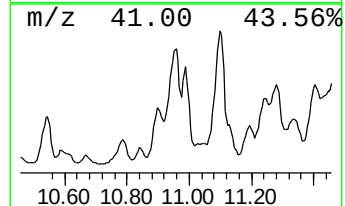
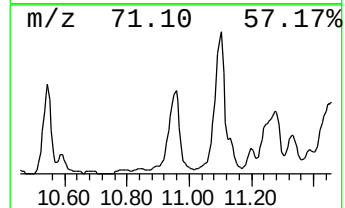
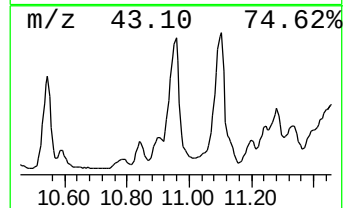
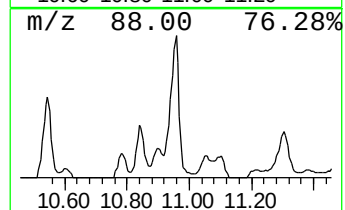
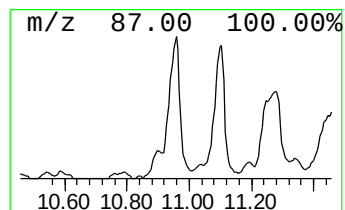
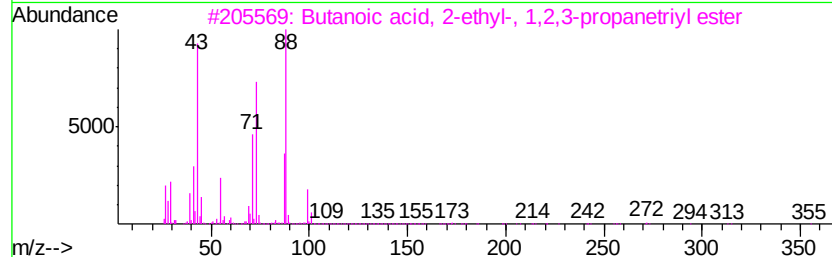
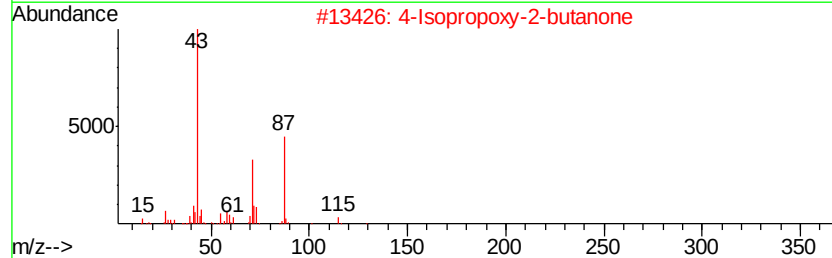
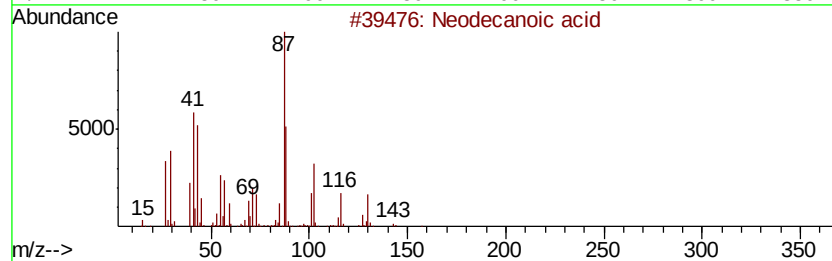
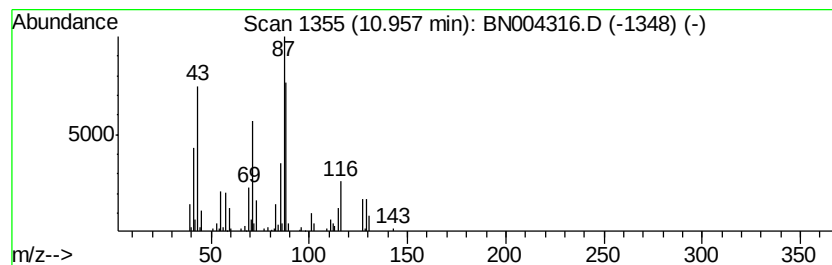
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-04 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.96	9.83 ng/ul	193144	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Neodecanoic acid	172	C10H20O2	026896-20-8	32
2		4-Isopropoxy-2-butanone	130	C7H14O2	032541-58-5	25
3		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	25
4		Methyl 2,3,4-tri-O-methyl-6-deox...	220	C10H20O5	072983-16-5	22
5		Heptanoic acid, 2,4-dimethyl-, m...	172	C10H20O2	018524-86-2	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
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Sample : J6593-01
Misc :
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Instrument :
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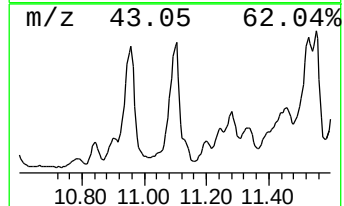
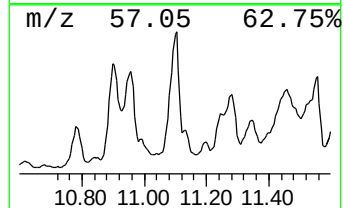
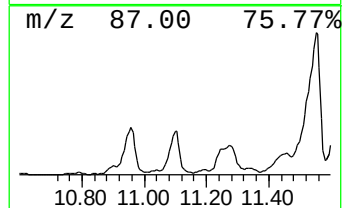
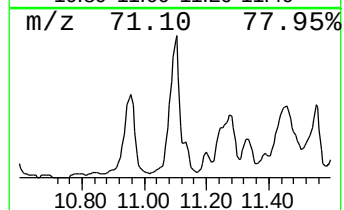
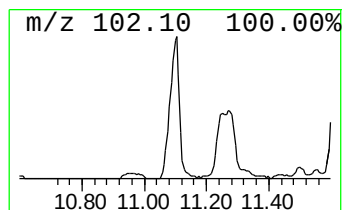
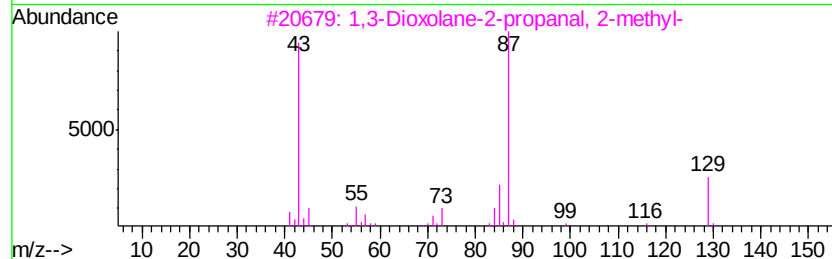
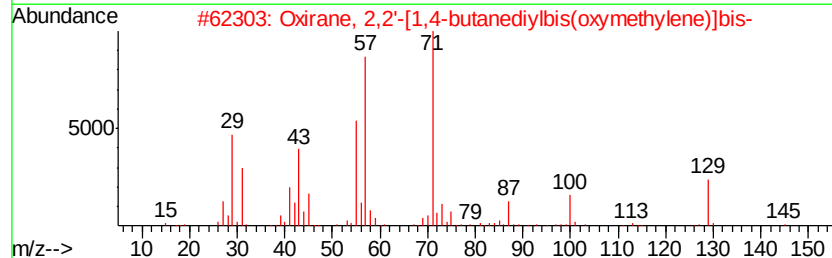
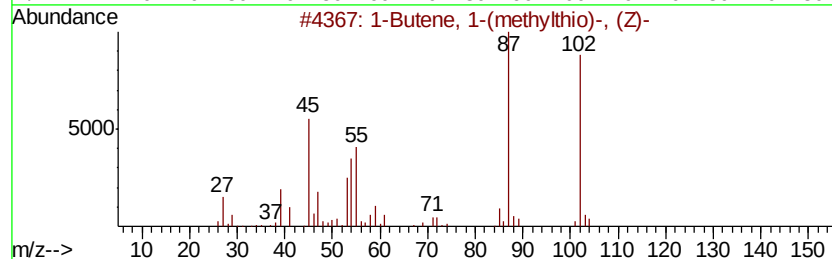
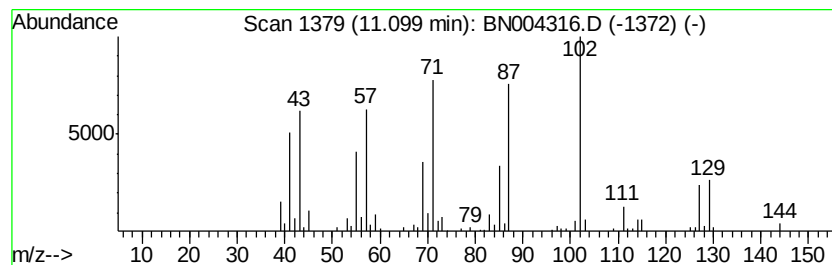
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	21.59 ng/ul	424088	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butene, 1-(methylthio)-, (Z)-	102	C5H10S	017414-15-2	38
2		Oxirane, 2,2'-[1,4-butanediylbis...	202	C10H18O4	002425-79-8	22
3		1,3-Dioxolane-2-propanal, 2-methyl-	144	C7H12O3	024108-29-0	22
4		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	14
5		2-Butene, 1-butoxy-3-methyl-	142	C9H18O	022094-02-6	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
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Misc :
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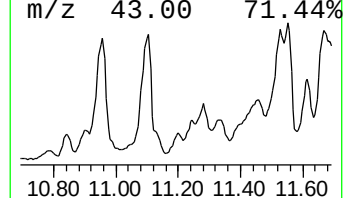
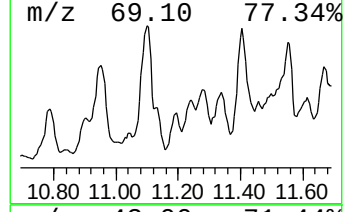
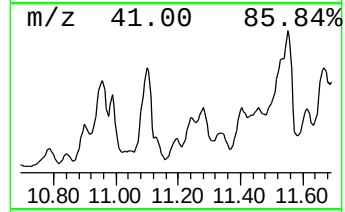
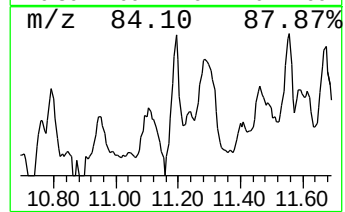
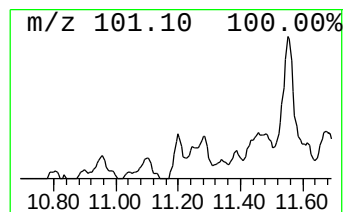
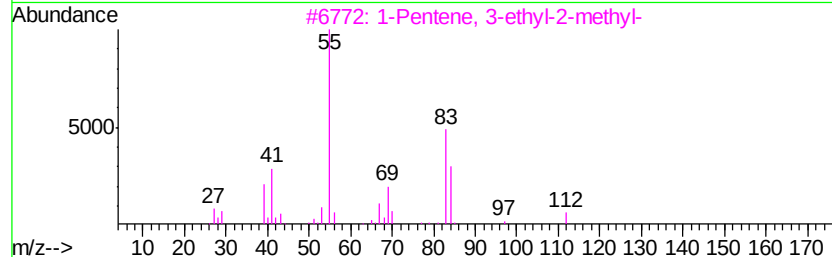
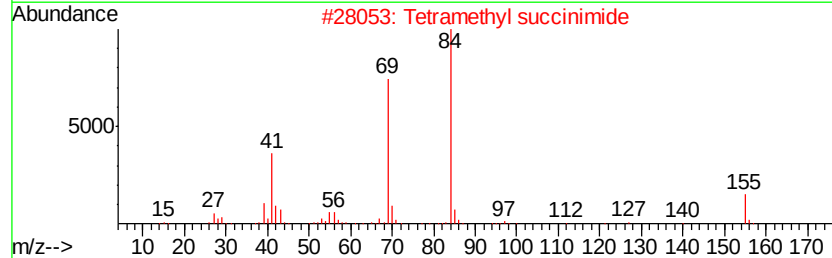
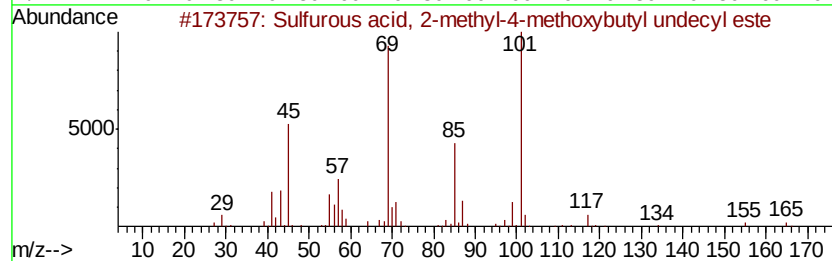
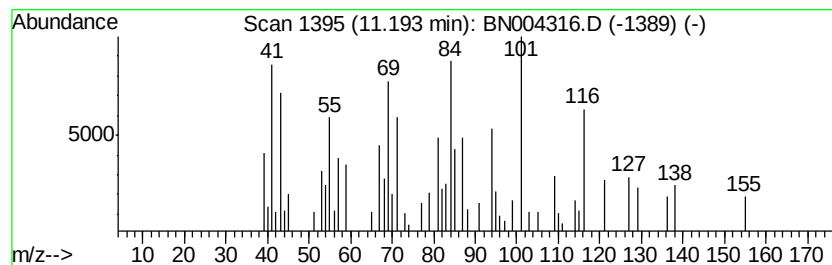
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-06 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.19	4.66 ng/ul	91521	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, 2-methyl-4-metho...	336	C17H36O4S	1000309-21-1	35
2		Tetramethyl succinimide	155	C8H13NO2	003566-61-8	22
3		1-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-66-6	22
4		3-Penten-2-one	84	C5H8O	000625-33-2	22
5		2H-Thiopyran, tetrahydro-2-methyl-	116	C6H12S	005161-16-0	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
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Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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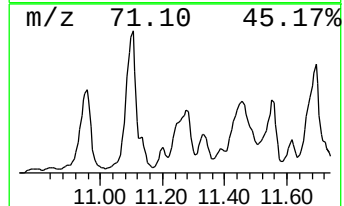
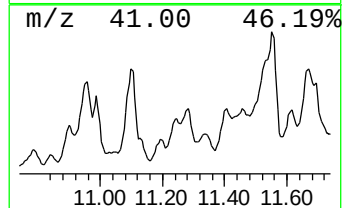
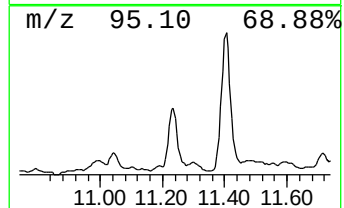
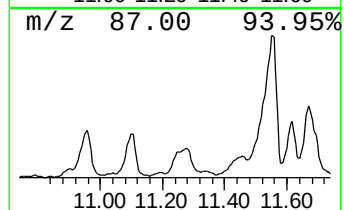
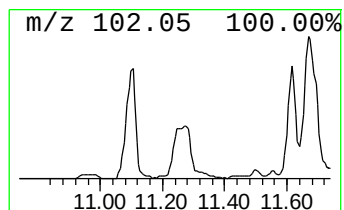
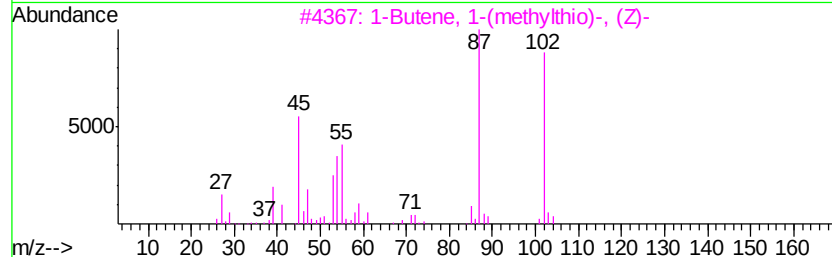
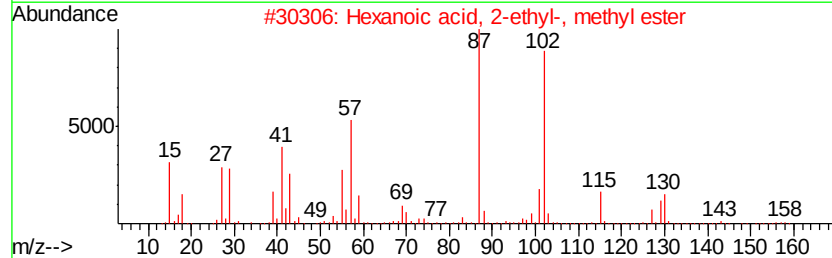
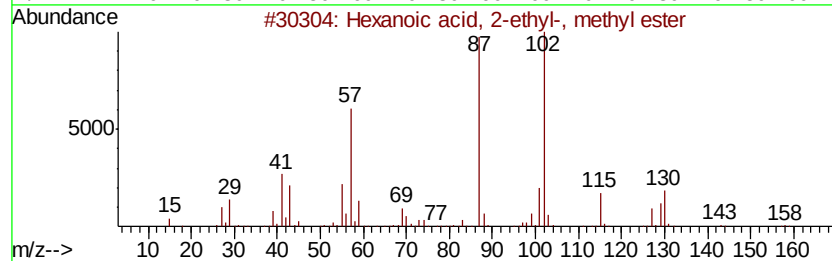
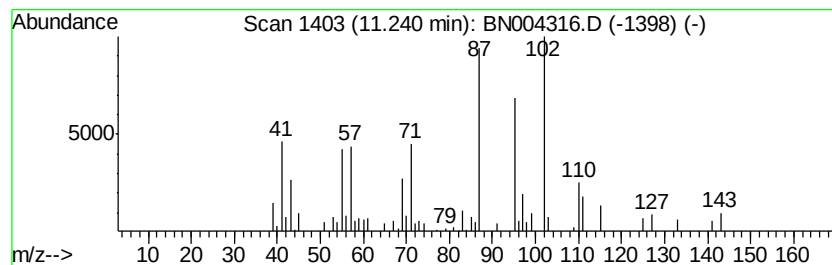
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-07 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.24	5.56 ng/ul	109327	Napthalene-d8	10.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-, methyl ...	158	C9H18O2	000816-19-3	38
2			Hexanoic acid, 2-ethyl-, methyl ...	158	C9H18O2	000816-19-3	38
3			1-Butene, 1-(methylthio)-, (Z)-	102	C5H10S	017414-15-2	38
4			Hexanoic acid, 2-ethyl-, methyl ...	158	C9H18O2	000816-19-3	32
5			1-Butene, 1-(methylthio)-, (E)-	102	C5H10S	017414-27-6	32



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
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Misc :
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ClientSampled :
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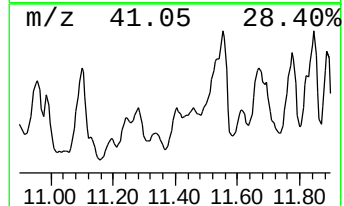
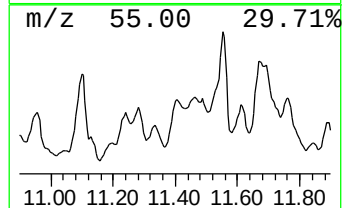
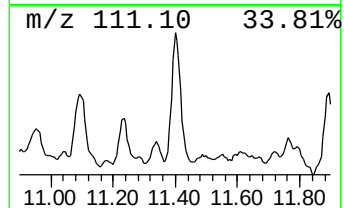
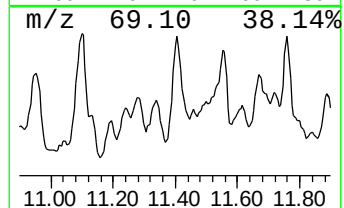
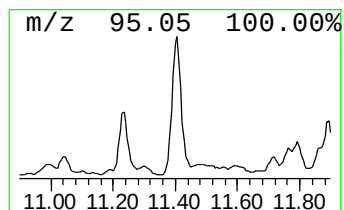
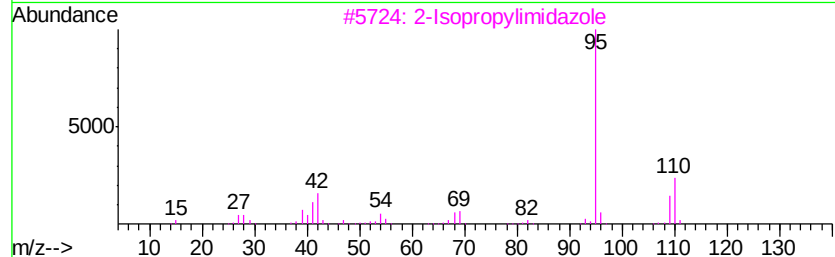
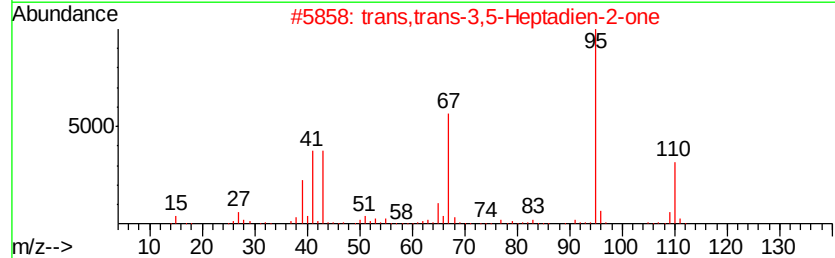
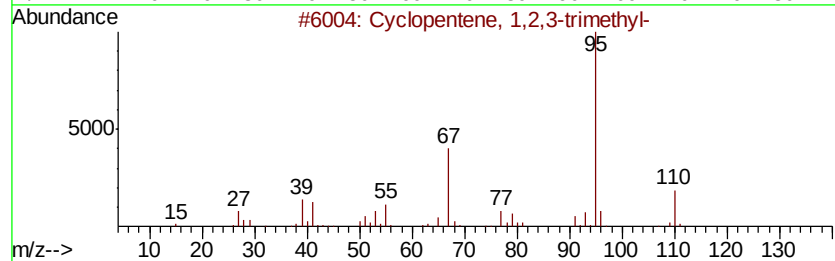
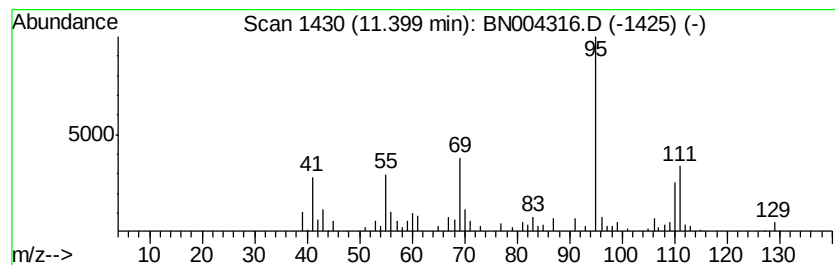
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Cyclopentene, 1,2,3-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	10.50 ng/ul	206371	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	50
2		trans,trans-3,5-Heptadien-2-one	110	C7H10O	018402-90-9	50
3		2-Isopropylimidazole	110	C6H10N2	036947-68-9	47
4		Ethanone, 1-(2-furanyl)-	110	C6H6O2	001192-62-7	47
5		5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BE803

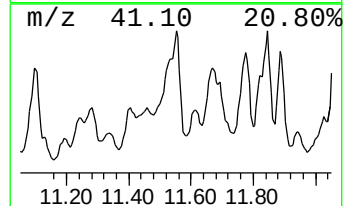
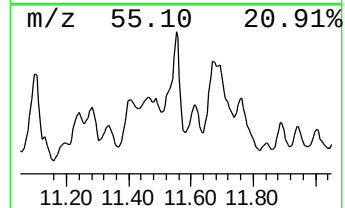
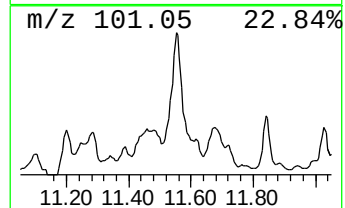
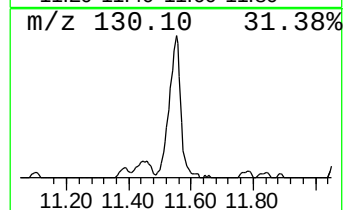
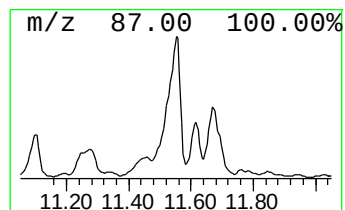
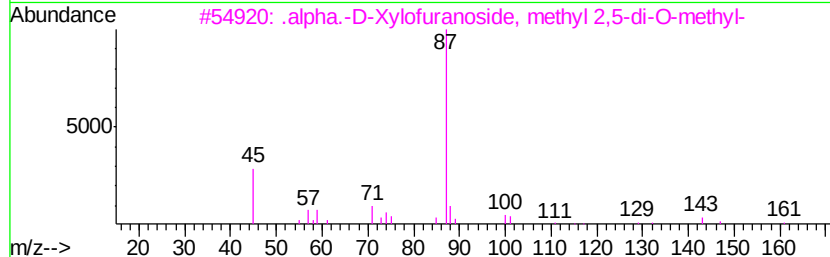
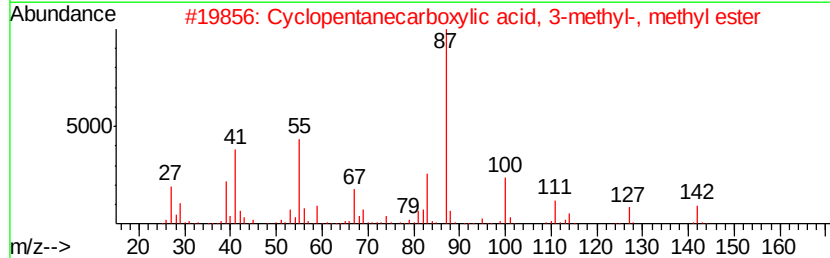
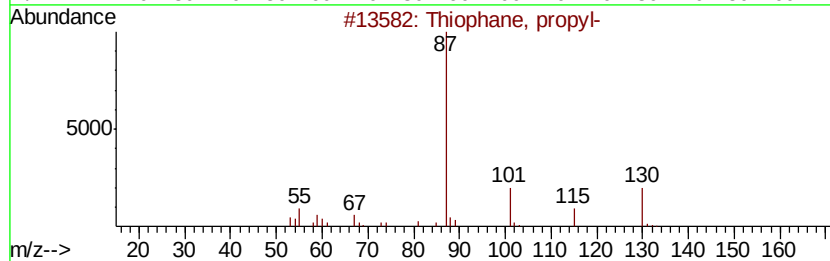
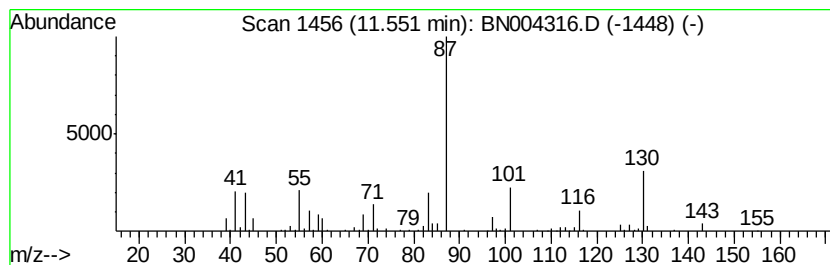
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Thiophane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	26.56 ng/ul	521745	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Thiophane, propyl-	130	C7H14S	001551-34-4	50
2		Cyclopentanecarboxylic acid, 3-m...	142	C8H14O2	024070-72-2	43
3		.alpha.-D-Xylofuranoside, methyl...	192	C8H16O5	035007-52-4	37
4		Heptanoic acid, 2-propyl-, methy...	186	C11H22O2	056247-53-1	37
5		1,3-Dioxolane, 2-butyl-4-methyl-	144	C8H16O2	074094-60-3	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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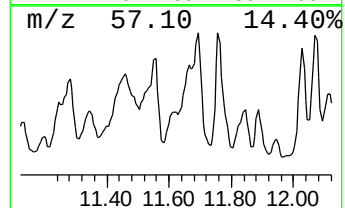
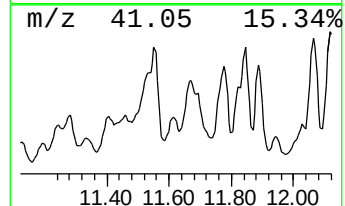
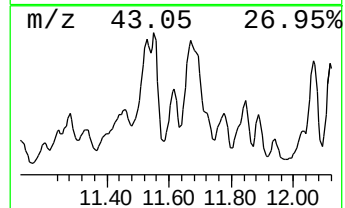
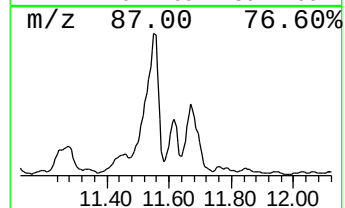
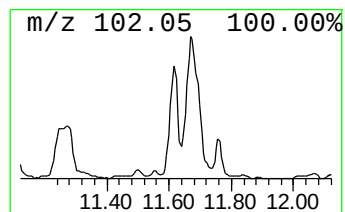
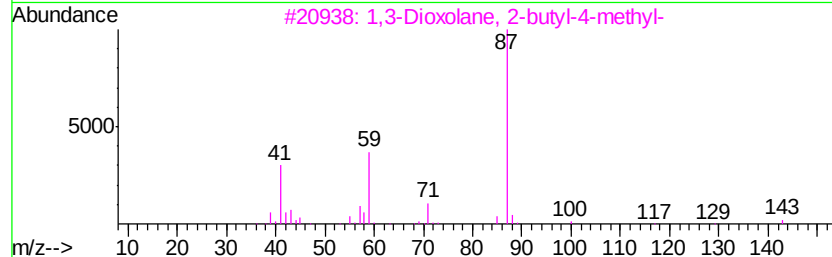
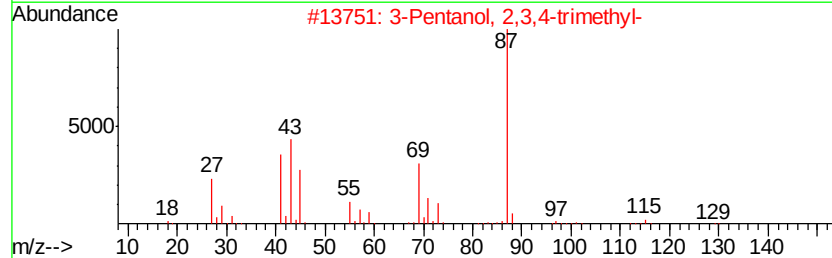
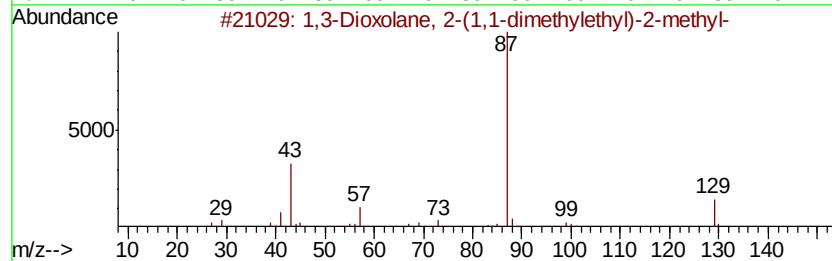
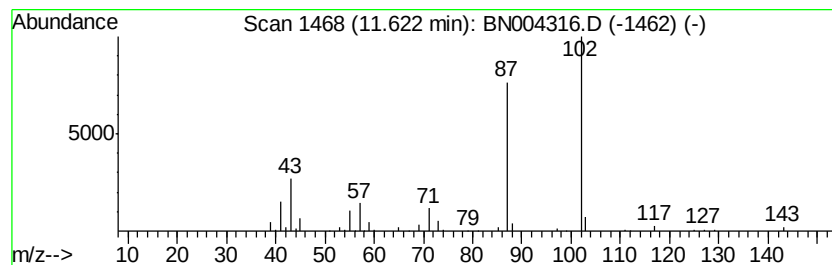
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-08 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	4.55 ng/ul	89455	Napthalene-d8	10.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 2-(1,1-dimethylet...	144	C8H16O2	006135-54-2	25
2			3-Pentanol, 2,3,4-trimethyl-	130	C8H18O	003054-92-0	23
3			1,3-Dioxolane, 2-butyl-4-methyl-	144	C8H16O2	074094-60-3	23
4			1,3-Dioxolane, 2-methyl-2-(2-met...	144	C8H16O2	002035-08-7	17
5			1,3-Dioxane	88	C4H8O2	000505-22-6	17



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
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Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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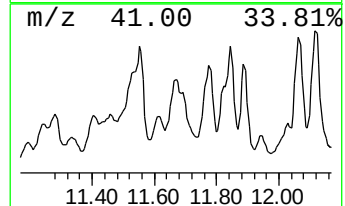
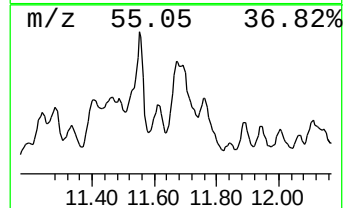
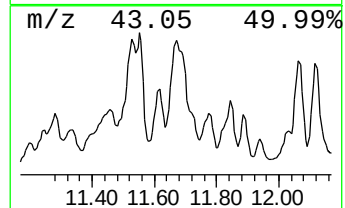
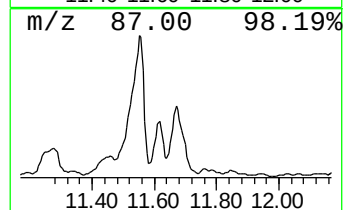
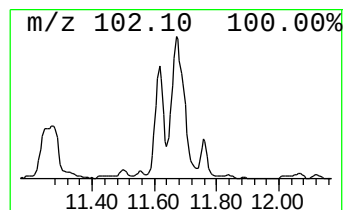
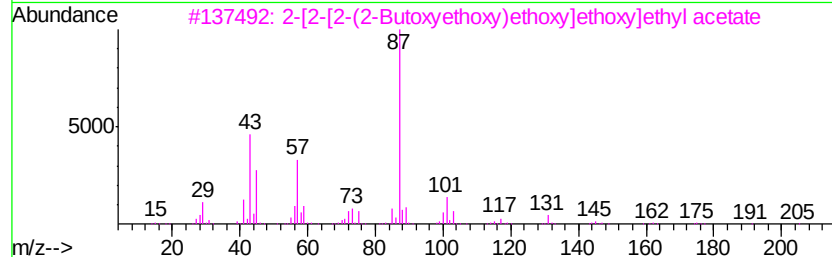
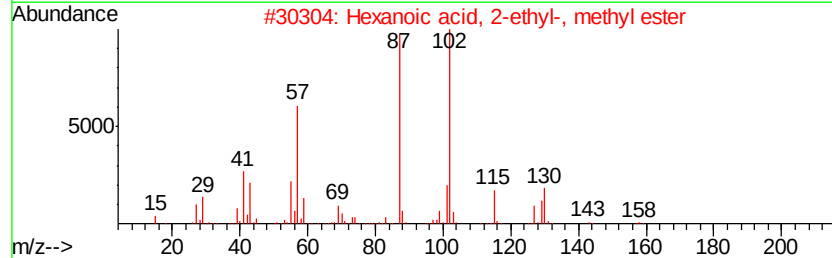
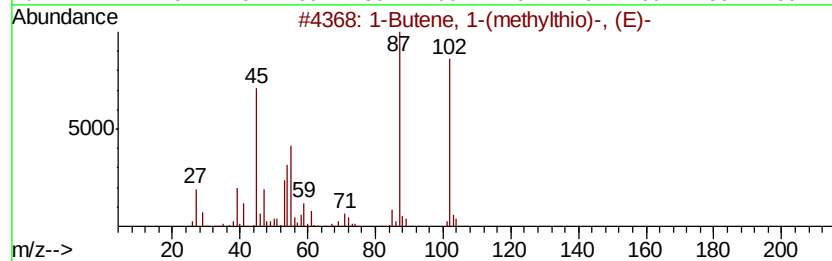
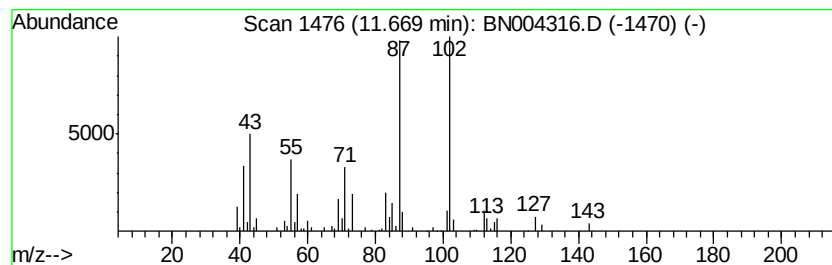
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1-Butene, 1-(methylthio)-, ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	12.55 ng/ul	246543	Napthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butene, 1-(methylthio)-, (E)-	102	C5H10S	017414-27-6	50
2		Hexanoic acid, 2-ethyl-, methyl ...	158	C9H18O2	000816-19-3	50
3		2-[2-[2-(2-Butoxyethoxy)ethoxy]e...	292	C14H28O6	1000351-93-9	35
4		Hexanoic acid, 2-ethyl-, methyl ...	158	C9H18O2	000816-19-3	33
5		Pentanoic acid, 2-ethyl-, methyl...	144	C8H16O2	000816-16-0	33



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
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Sample : J6593-01
Misc :
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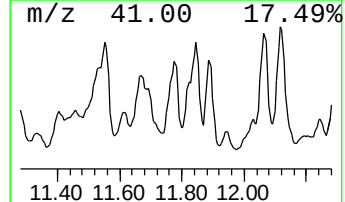
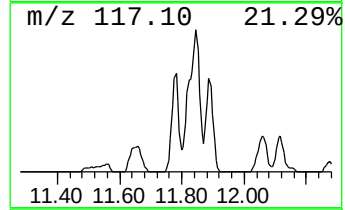
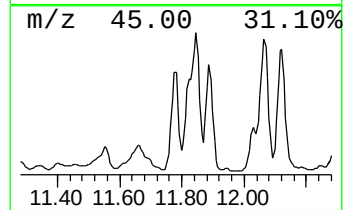
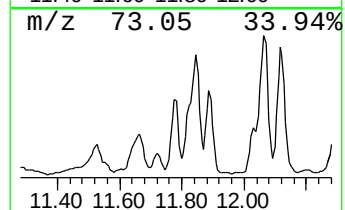
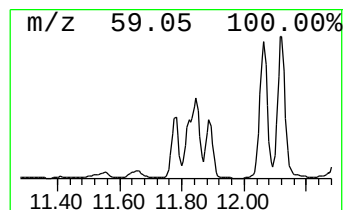
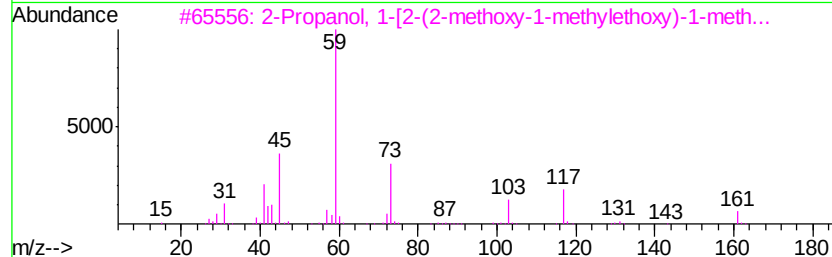
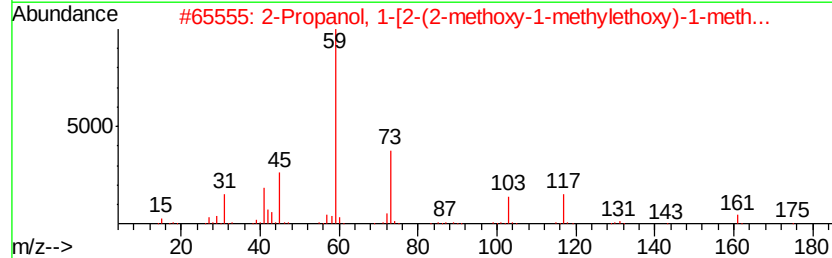
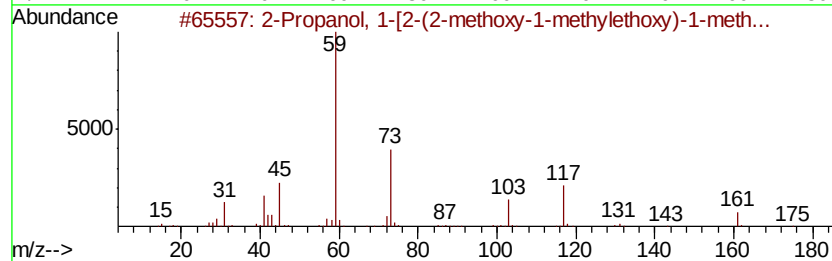
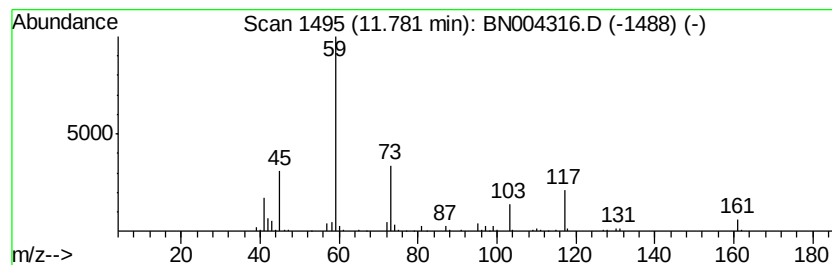
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	14.06 ng/ul	276130	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	86
2		2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	86
3		2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	78
4		2,5,8,11-Tetraoxatetradecan-13-o...	264	C13H28O5	020324-34-9	72
5		2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
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Misc :
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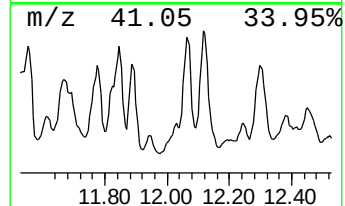
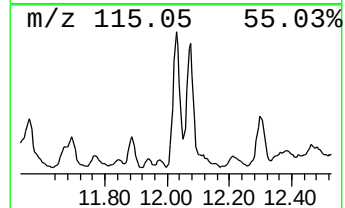
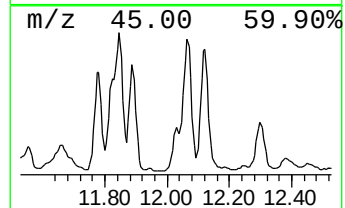
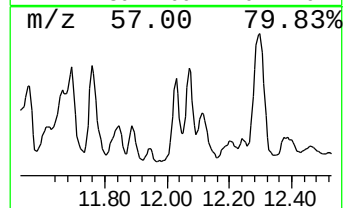
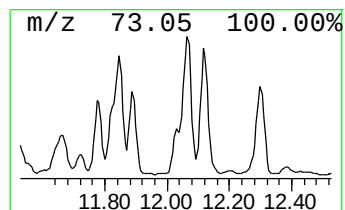
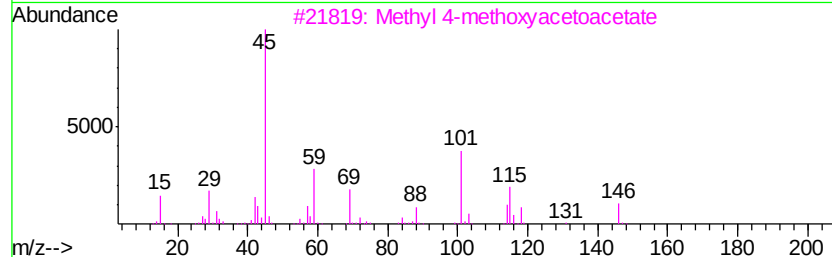
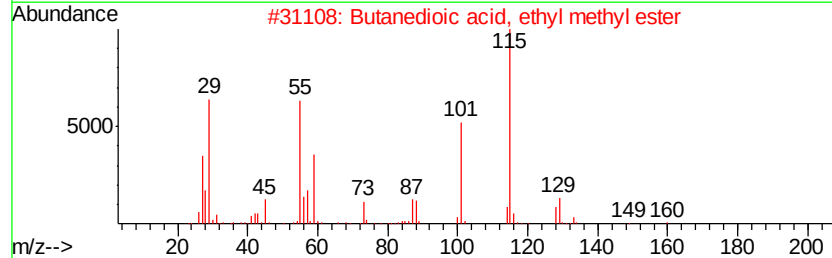
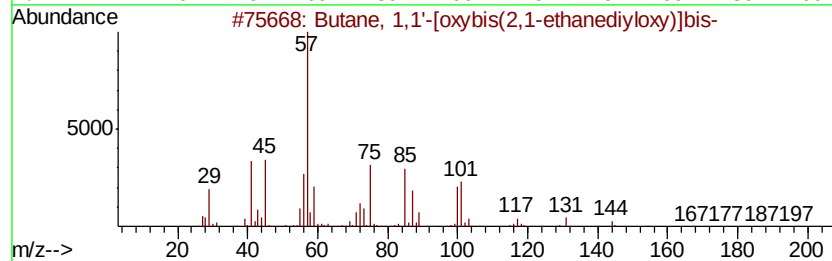
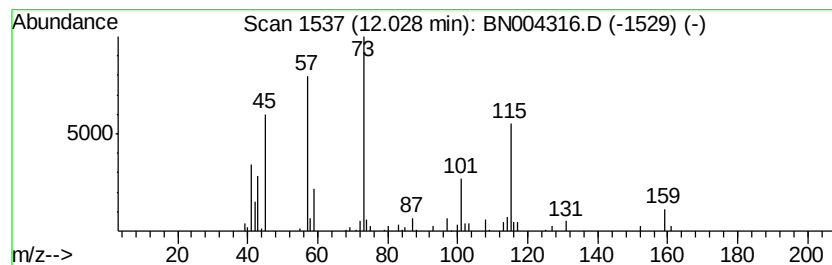
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-09 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	5.96 ng/ul	117081	Naphthalene-d8	10.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Butane, 1,1'-[oxybis(2,1-ethaned...	218 C12H26O3	000112-73-2 43
2		Butanedioic acid, ethyl methyl e...	160 C7H12O4	000627-73-6 43
3		Methyl 4-methoxyacetoacetate	146 C6H10O4	041051-15-4 37
4		2,13-Dibora-1,3,6,9,12,14,17,20-...	376 C16H34B2O8	1000150-01-2 32
5		Tripropylsilane	158 C9H22Si	000998-29-8 28



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
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Acq On : 03 Jan 2019 13:58
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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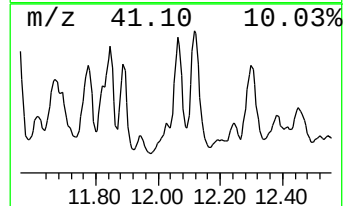
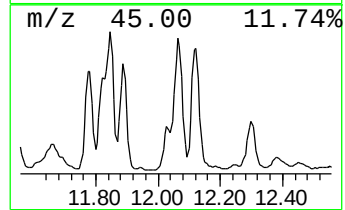
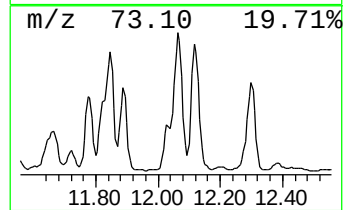
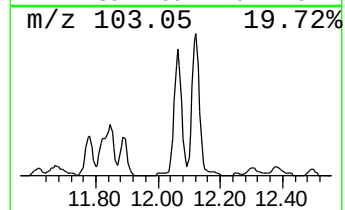
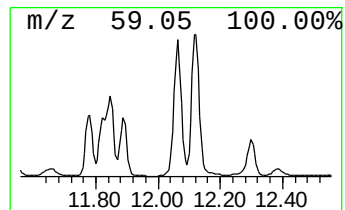
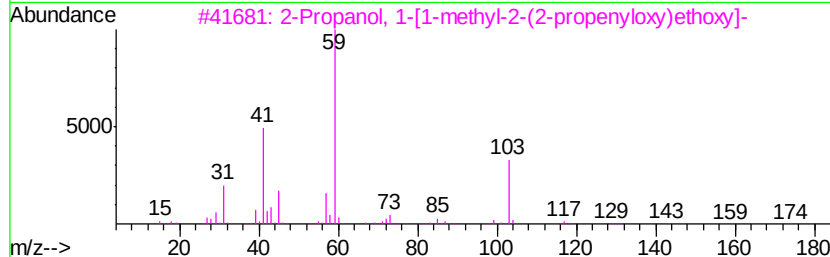
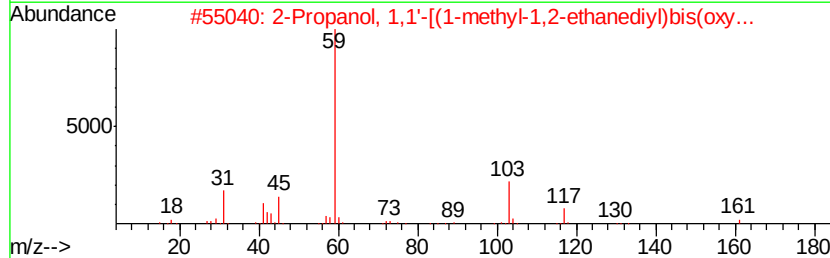
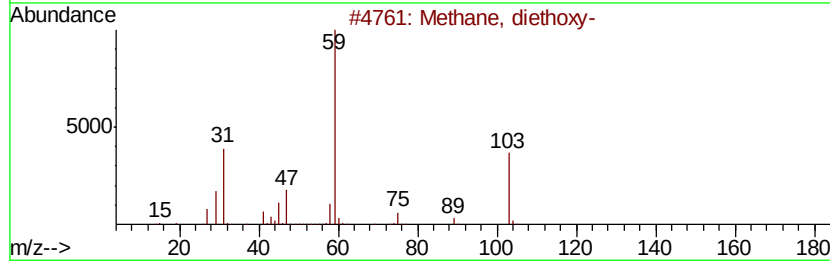
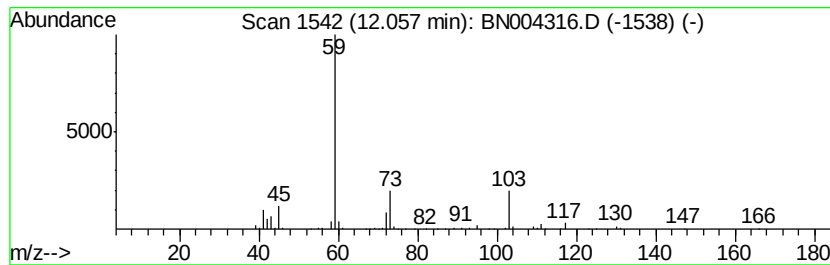
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Methane, diethoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.06	20.02 ng/ul	393252	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methane, diethoxy-	104	C5H12O2	000462-95-3	64
2		2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	64
3		2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	64
4		2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	59
5		Methane, diethoxy-	104	C5H12O2	000462-95-3	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
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Acq On : 03 Jan 2019 13:58
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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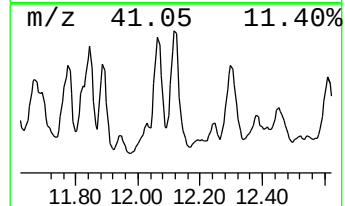
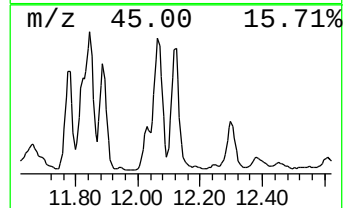
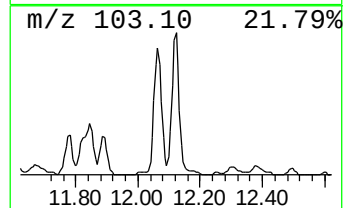
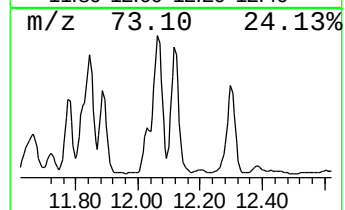
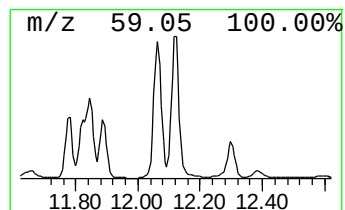
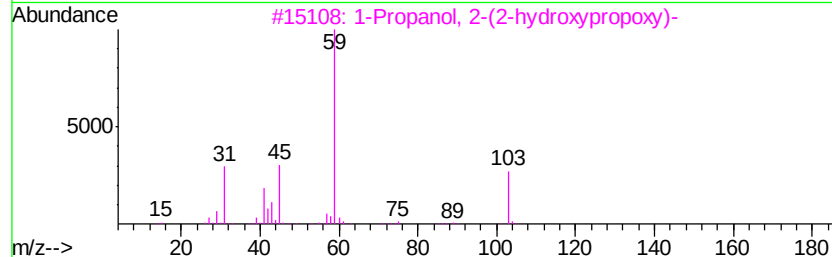
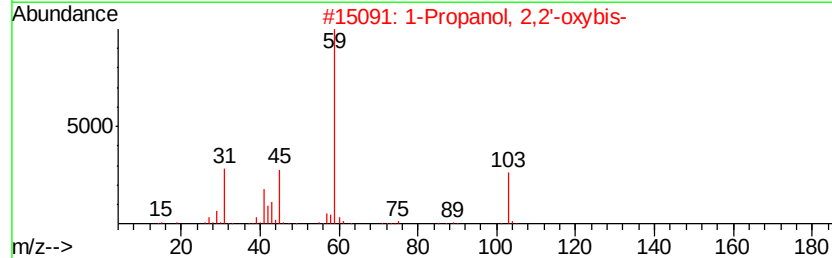
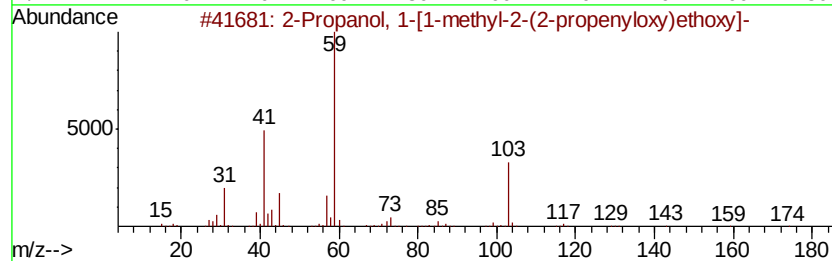
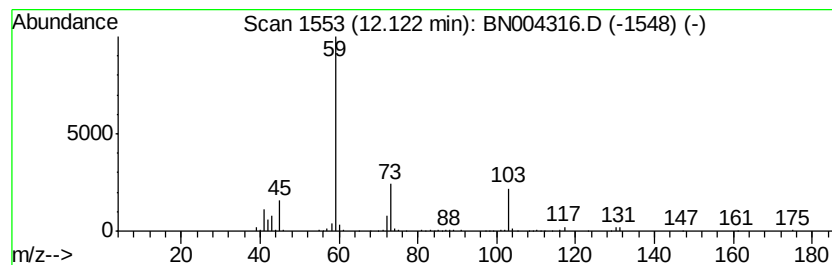
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 2-Propanol, 1-[1-methyl-2-(... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	27.63 ng/ul	542756	Napthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	64
2		1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	59
3		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	53
4		Tetraglyme	222	C10H22O5	000143-24-8	50
5		Propane, 1,2-dimethoxy-	104	C5H12O2	007778-85-0	50



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Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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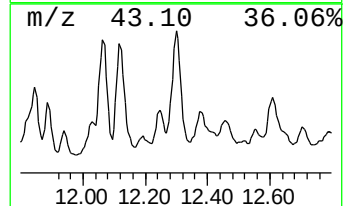
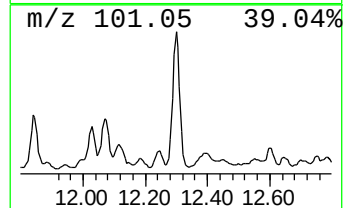
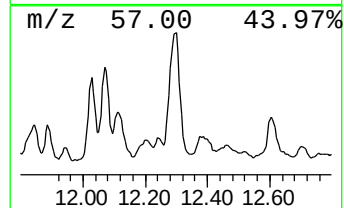
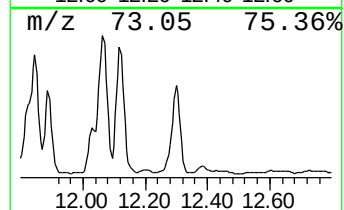
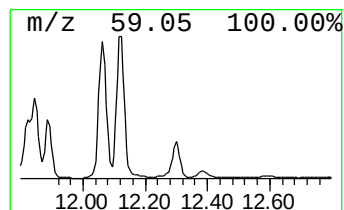
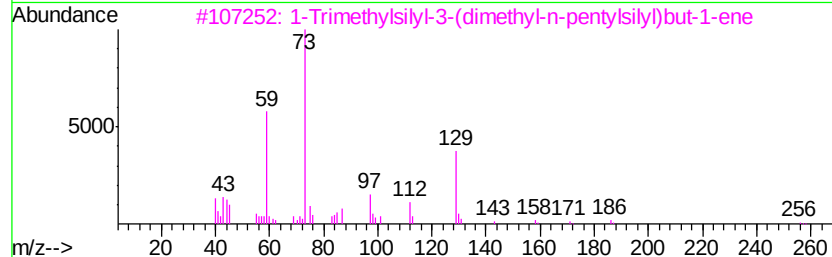
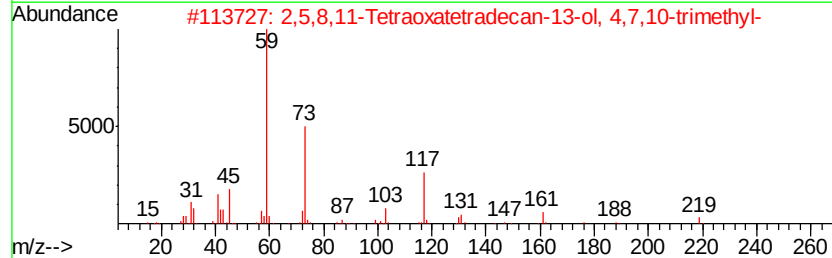
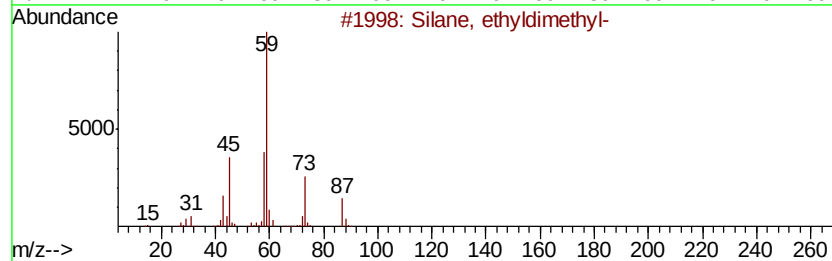
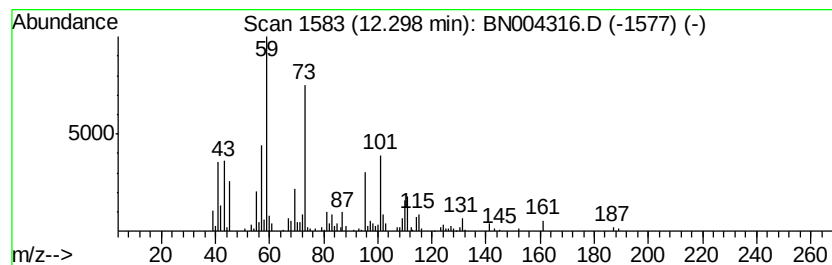
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 unknown-10 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	20.66 ng/ul	405982	Napthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	46
2		2,5,8,11-Tetraoxatetradecan-13-o...	264	C13H28O5	020324-34-9	43
3		1-Trimethylsilyl-3-(dimethyl-n-p...	256	C14H32Si2	136935-52-9	43
4		Silane, butyltrimethyl-	130	C7H18Si	001000-49-3	38
5		Trimethyl(n-pentyl)silane	144	C8H20Si	001641-49-2	37



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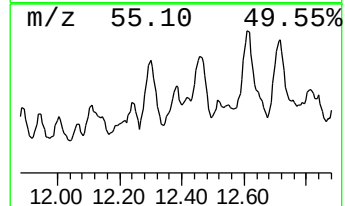
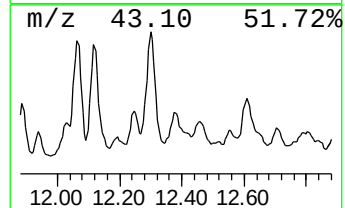
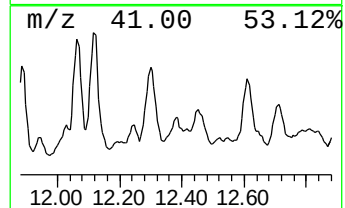
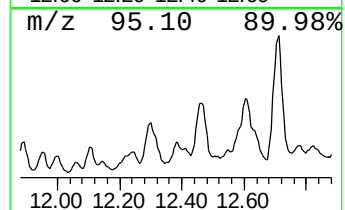
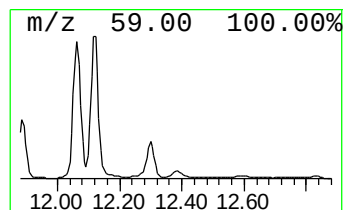
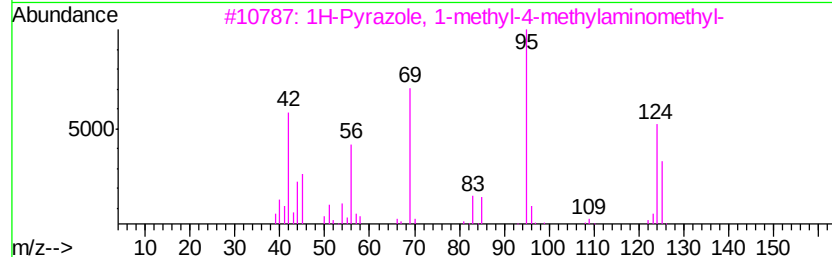
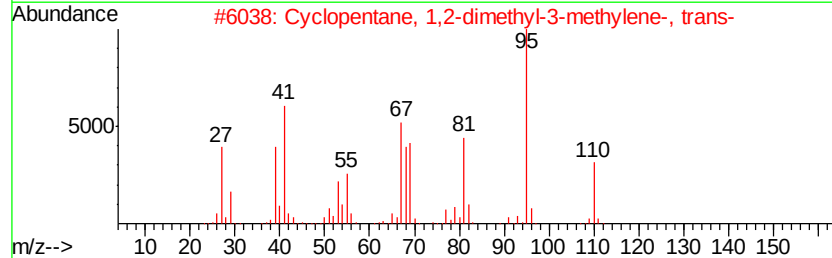
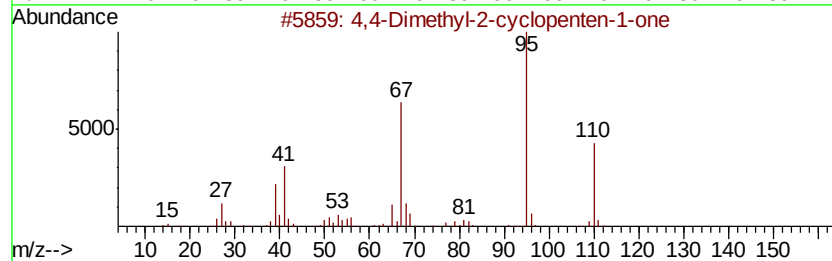
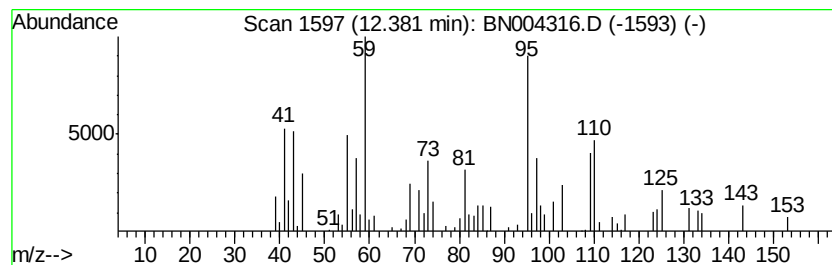
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 unknown-11 Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.38	3.68 ng/ul	72265	Naphthalene-d8	10.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,4-Dimethyl-2-cyclopenten-1-one	110	C7H10O	022748-16-9	22
2	Cyclopentane, 1,2-dimethyl-3-met...	110	C8H14	1000150-99-3	14
3	1H-Pyrazole, 1-methyl-4-methylam...	125	C6H11N3	1000294-10-8	14
4	Ethanone, 1-(2-furanyl)-	110	C6H6O2	001192-62-7	14
5	2,4-Hexadiene, 2,5-dimethyl-	110	C8H14	000764-13-6	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
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Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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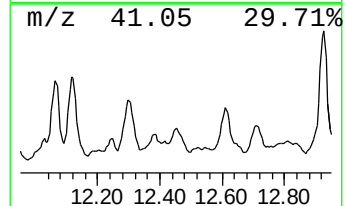
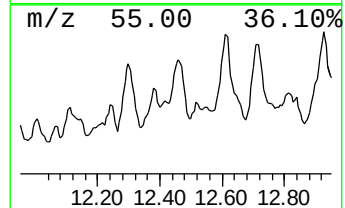
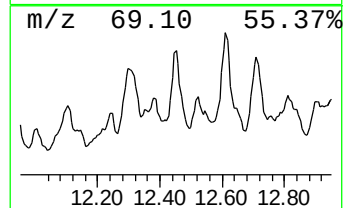
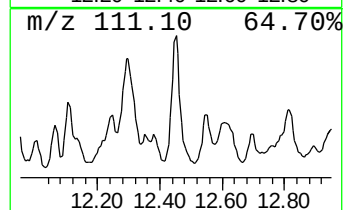
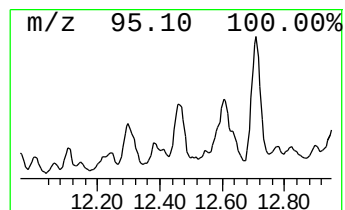
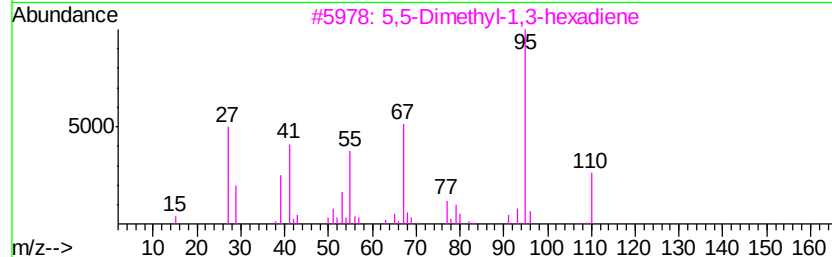
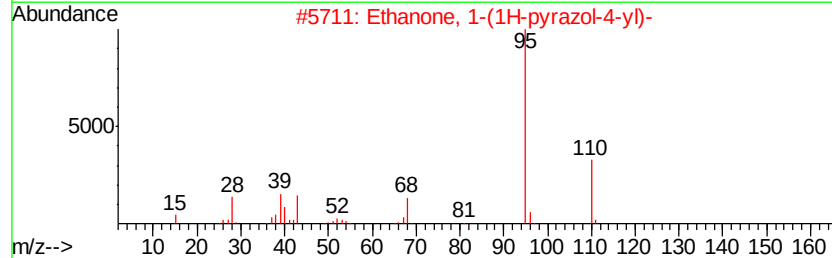
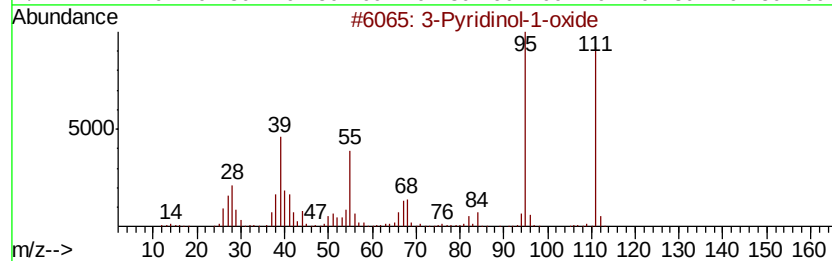
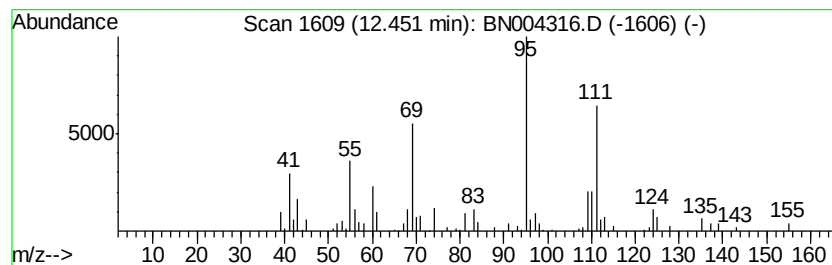
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 unknown-12 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	8.24 ng/ul	161856	Naphthalene-d8	10.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pyridinol-1-oxide	111	C5H5NO2	006602-28-4	46
2			Ethanone, 1-(1H-pyrazol-4-yl)-	110	C5H6N2O	025016-16-4	35
3			5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	35
4			1,3-Dimethyl-5-n-decylcyclohexane	252	C18H36	086710-36-3	22
5			Cyclopentane, 1,1,3-trimethyl-3-...	166	C12H22	074421-09-3	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
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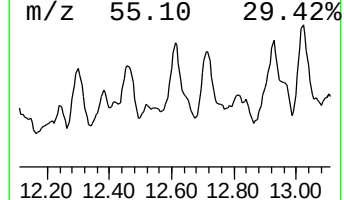
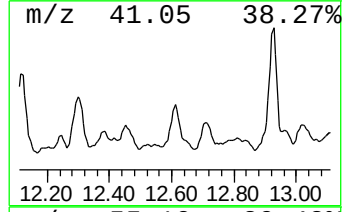
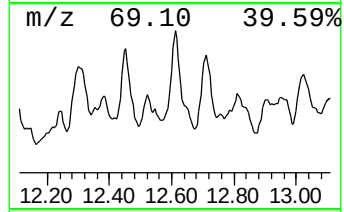
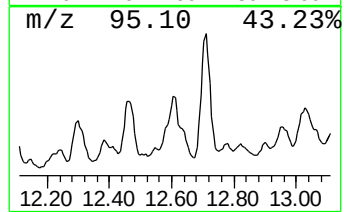
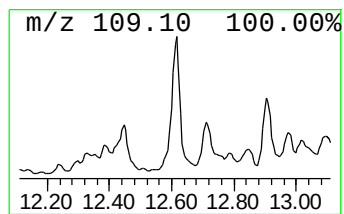
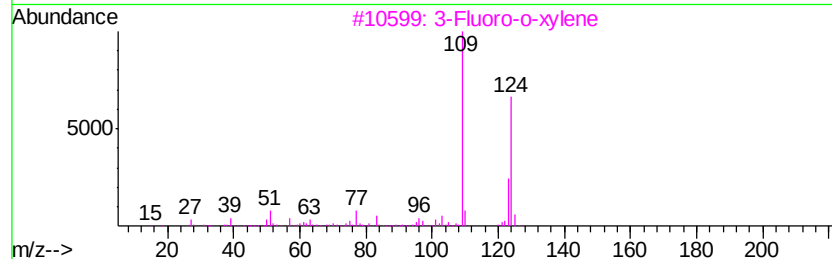
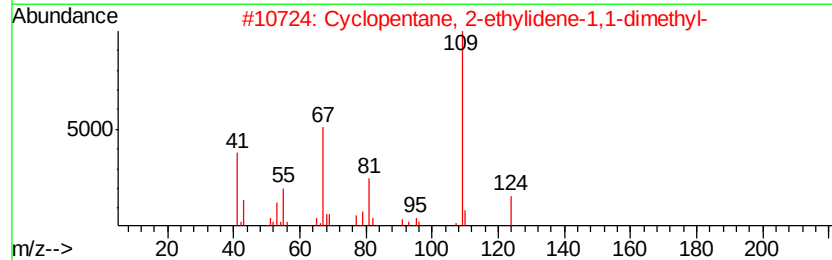
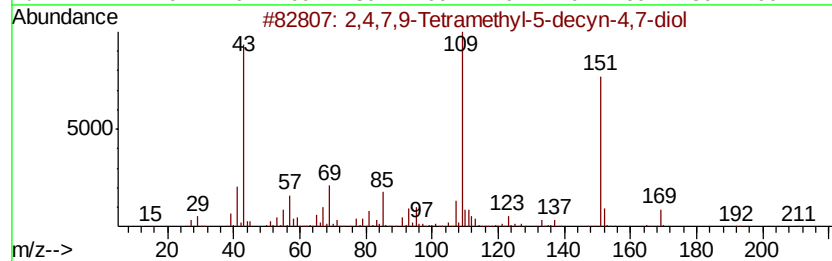
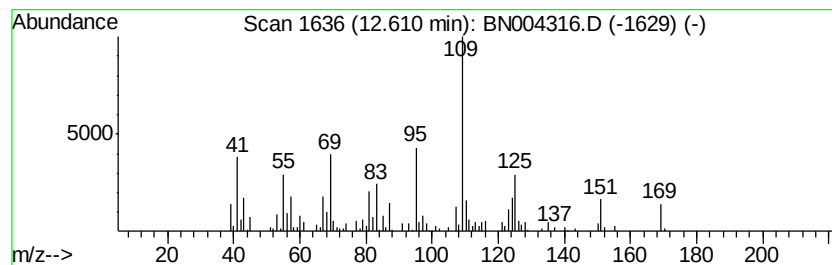
Instrument :
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 unknown-13 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	14.50 ng/ul	411960	Acenaphthene-d10	14.45
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226 C14H26O2	000126-86-3	43
2	Cyclopentane, 2-ethylidene-1,1-d...	124 C9H16	056324-66-4	38
3	3-Fluoro-o-xylene	124 C8H9F	000443-82-3	38
4	6,6-Dimethylhepta-2,4-diene	124 C9H16	1000195-03-3	38
5	1(3H)-Isobenzofuranone, 3a,4,5,7...	182 C10H14O3	054346-06-4	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
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Sample : J6593-01
Misc :
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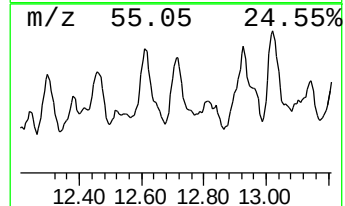
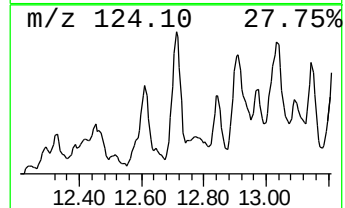
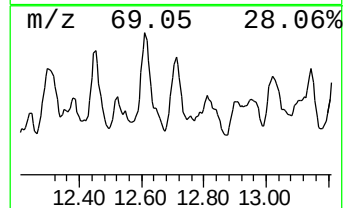
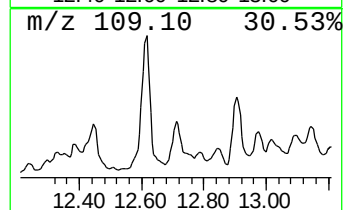
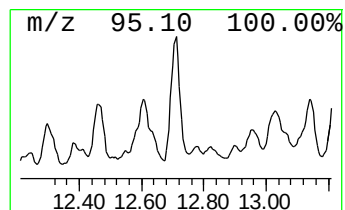
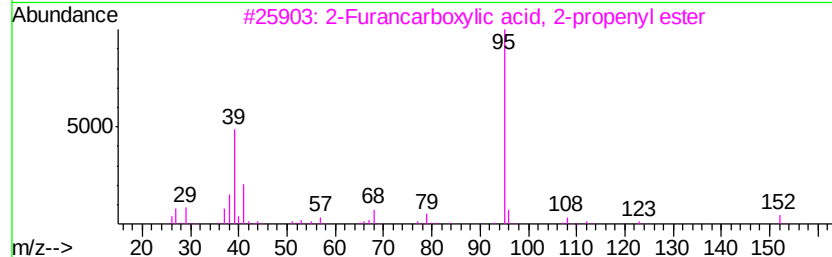
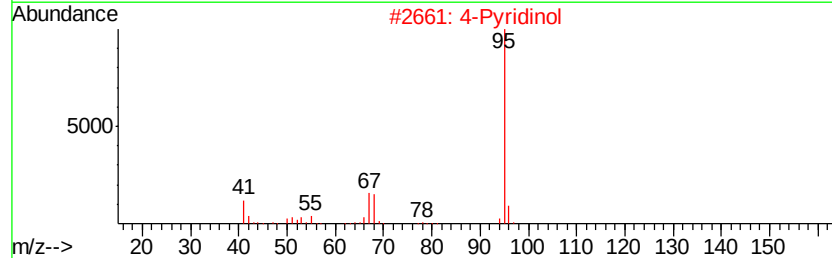
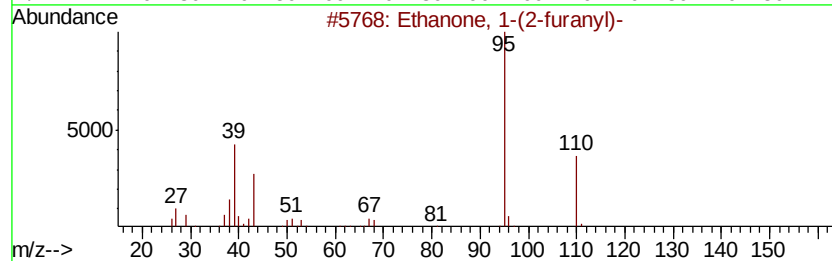
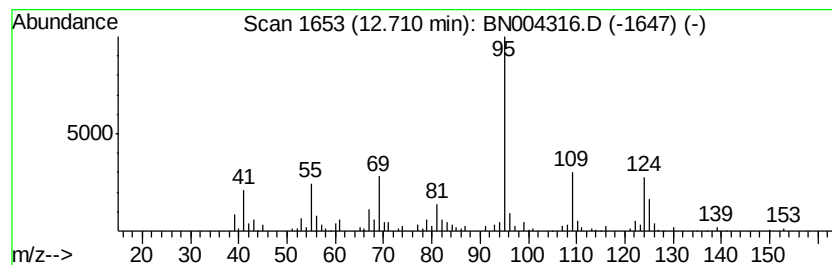
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-14 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	9.44 ng/ul	268267	Acenaphthene-d10	14.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 1-(2-furanyl)-	110	C6H6O2	001192-62-7	46
2		4-Pyridinol	95	C5H5NO	000626-64-2	46
3		2-Furancarboxylic acid, 2-propen...	152	C8H8O3	004208-49-5	43
4		1-Bromomethyl-2-chlorocyclohexane	210	C7H12BrCl	090009-23-7	43
5		3-Pyridinecarboxylic acid, 4-hyd...	139	C6H5NO3	000609-70-1	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
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ALS Vial : 4 Sample Multiplier: 1

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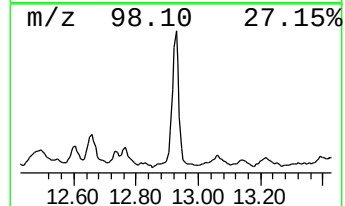
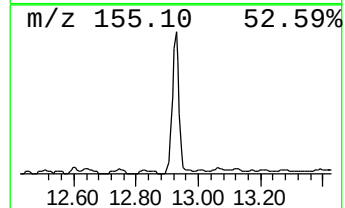
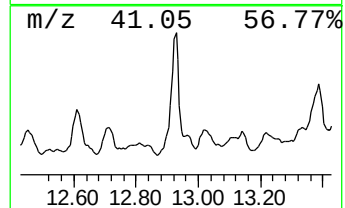
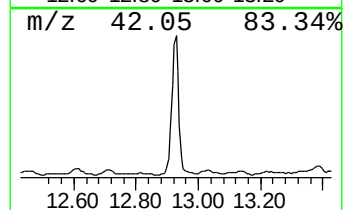
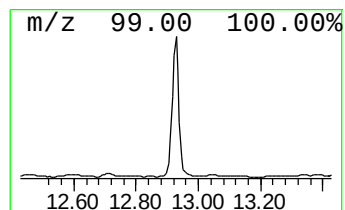
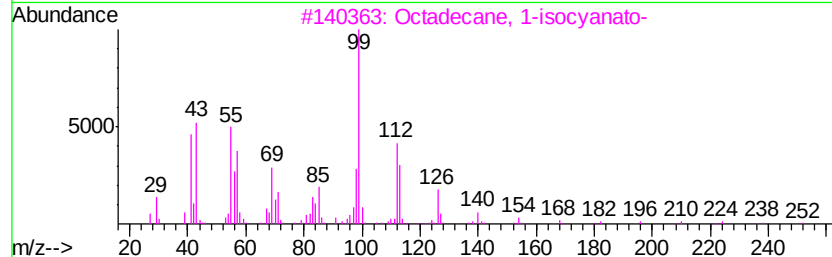
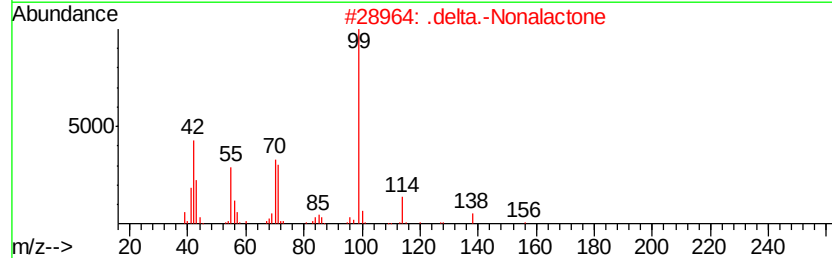
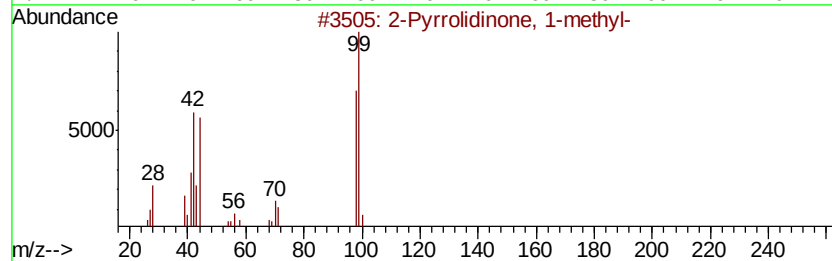
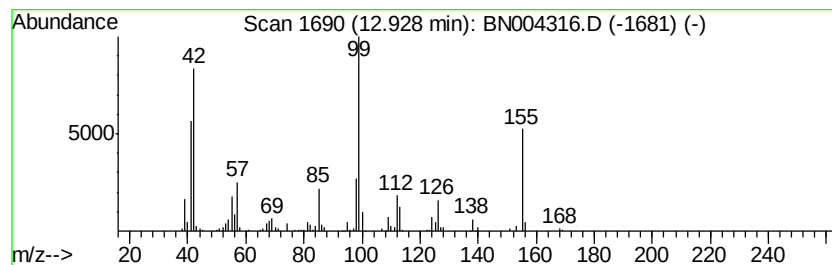
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-15 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	17.61 ng/ul	500544	Acenaphthene-d10	14.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pyrrolidinone, 1-methyl-	99	C5H9NO	000872-50-4	47
2		.delta.-Nonalactone	156	C9H16O2	003301-94-8	38
3		Octadecane, 1-isocyanato-	295	C19H37NO	000112-96-9	32
4		Oxazole, 4,5-dihydro-2,4,4-trime...	113	C6H11NO	001772-43-6	30
5		3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
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Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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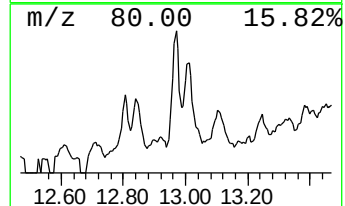
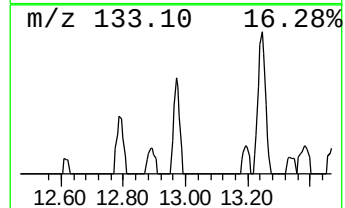
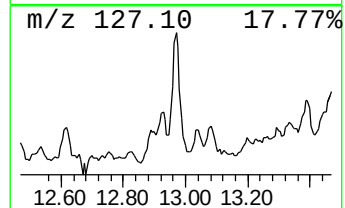
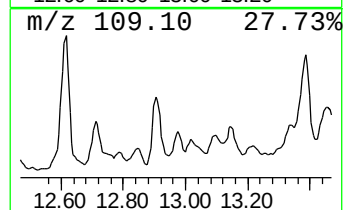
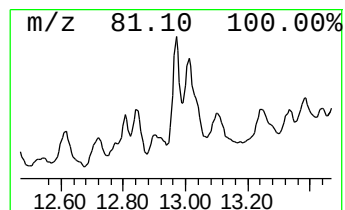
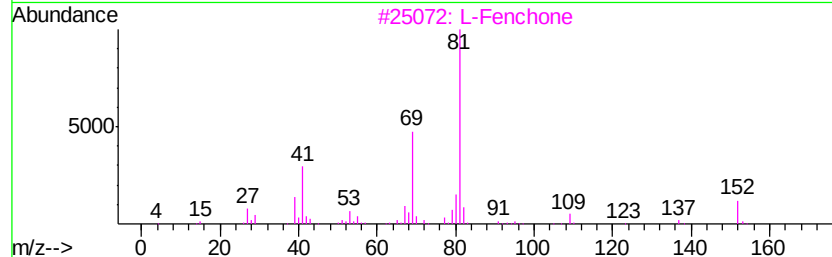
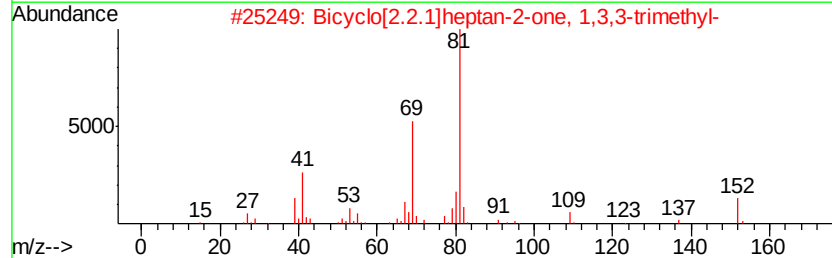
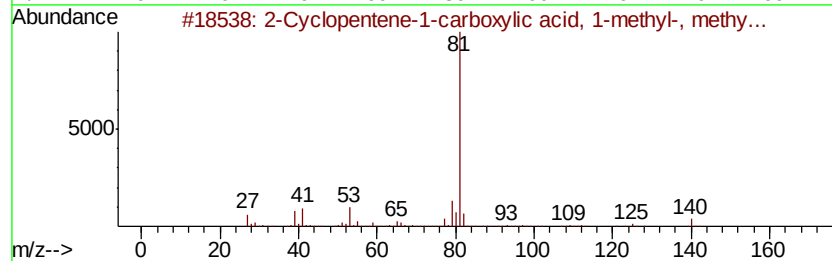
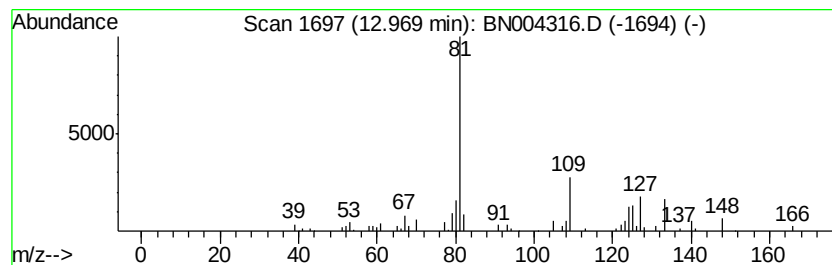
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 unknown-16 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	3.73 ng/ul	105912	Acenaphthene-d10	14.45

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclopentene-1-carboxylic acid...	140	C8H12O2	068317-73-7	46
2			Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	43
3			L-Fenchone	152	C10H16O	007787-20-4	37
4			2-n-Butyl furan	124	C8H12O	004466-24-4	37
5			2-Cyclopentene-1-acetaldehyde, 2...	166	C10H14O2	075332-42-2	33



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampled :
BE803

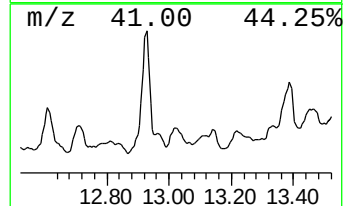
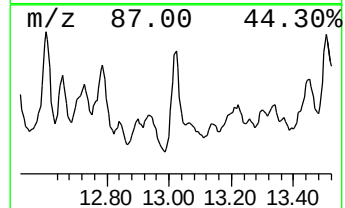
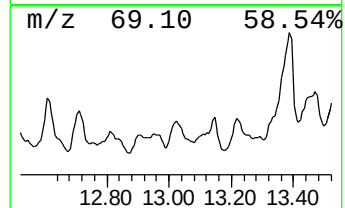
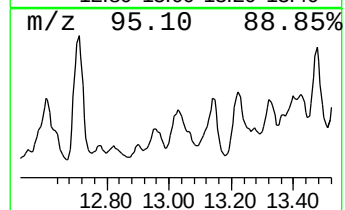
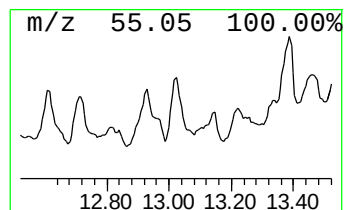
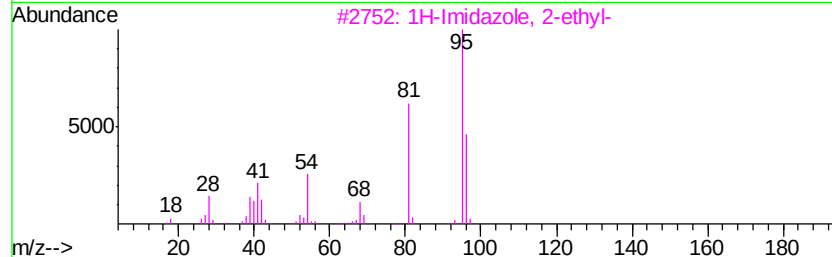
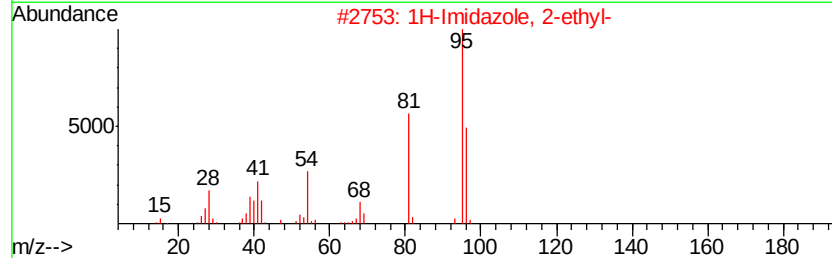
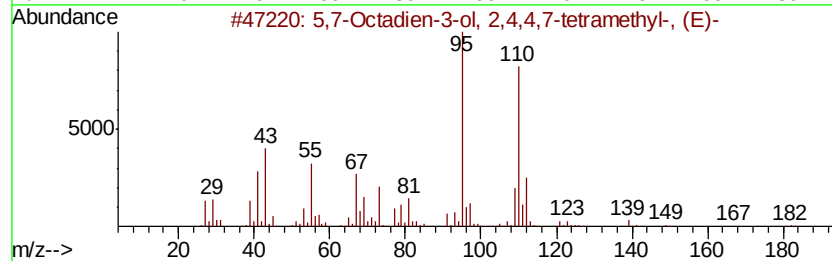
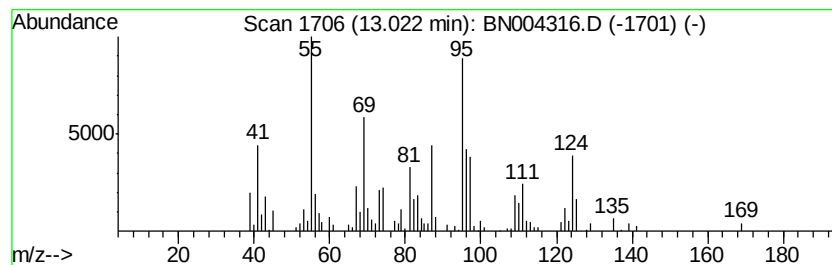
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 unknown-17 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	6.82 ng/ul	193957	Acenaphthene-d10	14.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5,7-Octadien-3-ol, 2,4,4,7-tetra...	182	C12H22O	077142-78-0	35
2		1H-Imidazole, 2-ethyl-	96	C5H8N2	001072-62-4	27
3		1H-Imidazole, 2-ethyl-	96	C5H8N2	001072-62-4	27
4		1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	25
5		1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

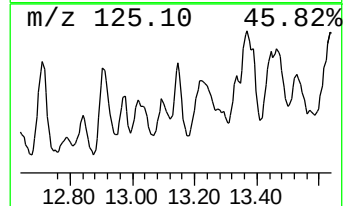
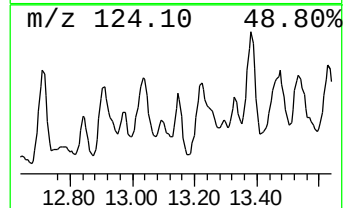
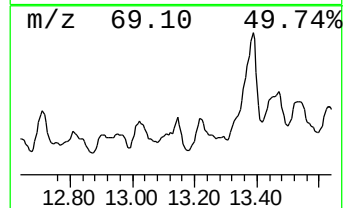
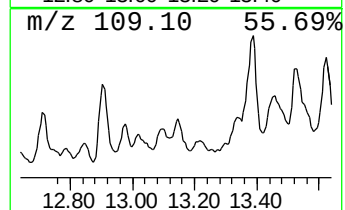
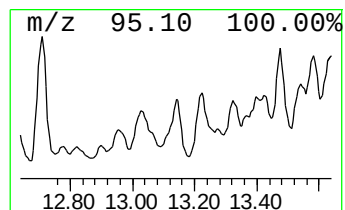
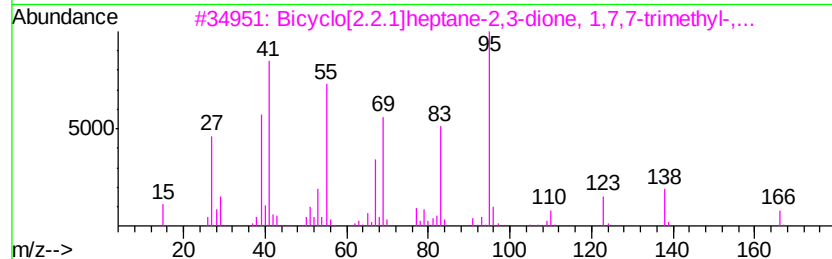
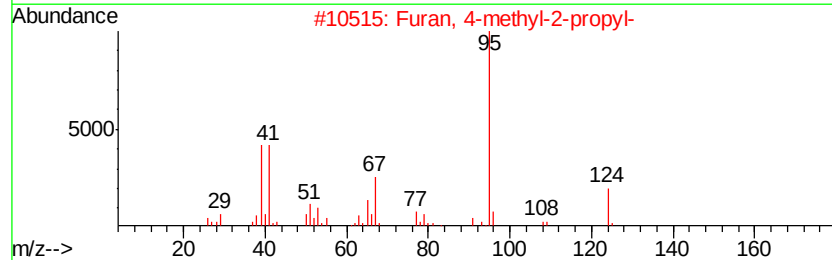
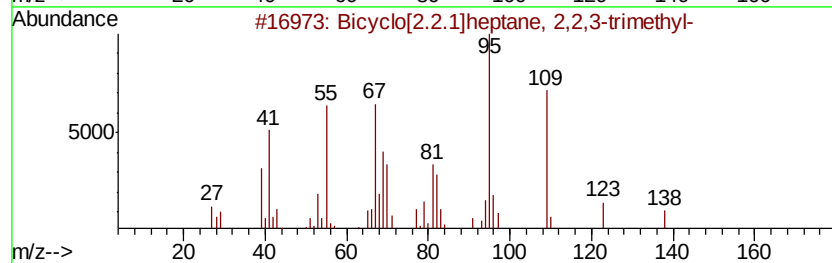
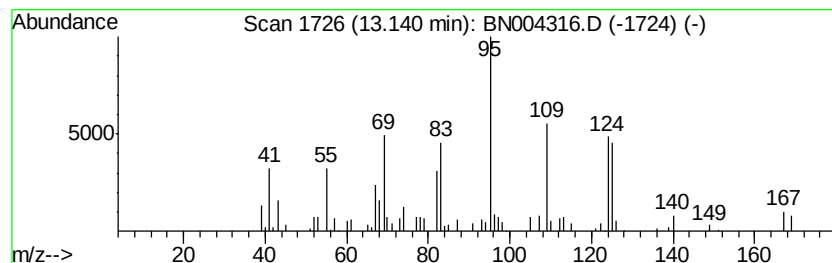
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 29 unknown-18 Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.14	3.53 ng/ul	100426	Acenaphthene-d10	14.45
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Bicyclo[2.2.1]heptane, 2,2,3-tri...	138 C10H18	000473-19-8	37
2	Furan, 4-methyl-2-propyl-	124 C8H12O	006148-37-4	35
3	Bicyclo[2.2.1]heptane-2,3-dione,...	166 C10H14O2	010334-26-6	35
4	5-Ethyl-2-furaldehyde	124 C7H8O2	023074-10-4	27
5	Silane, diethyldifluoro-	124 C4H10F2Si	000358-06-5	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

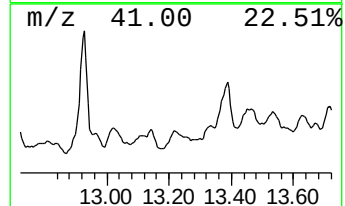
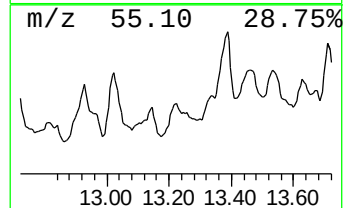
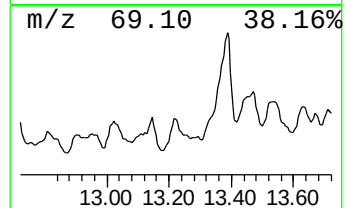
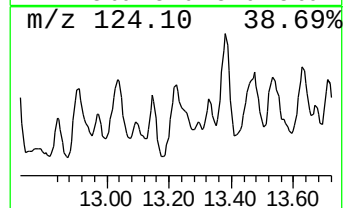
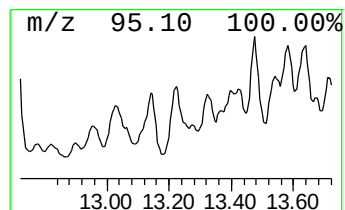
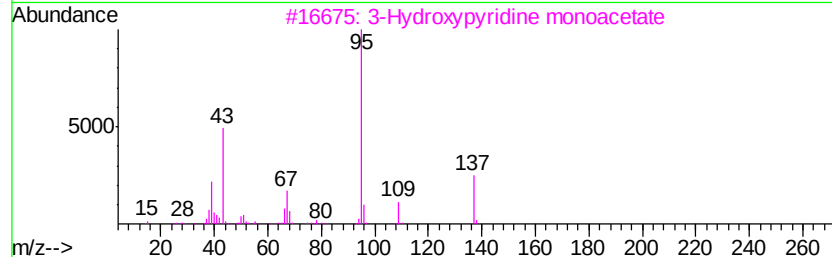
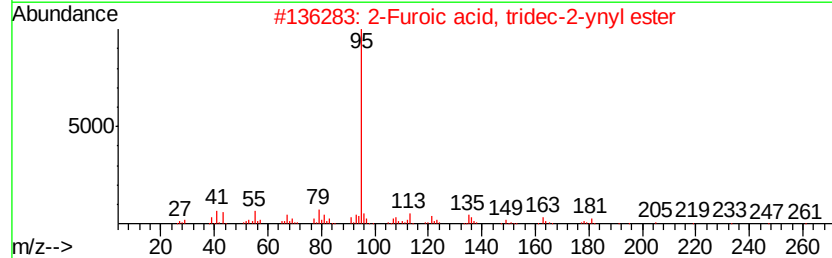
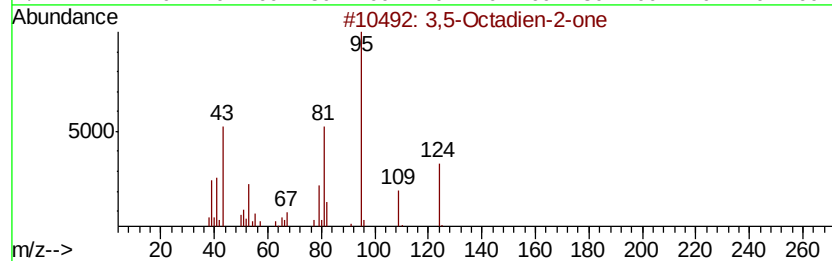
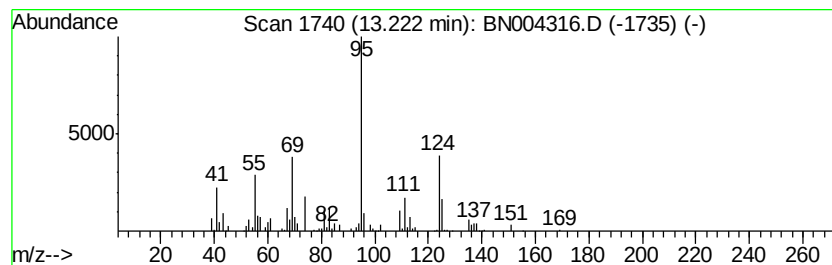
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 unknown-19 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	6.49 ng/ul	184565	Acenaphthene-d10	14.45
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	3,5-Octadien-2-one	124 C8H12O	038284-27-4	47
2	2-Furoic acid, tridec-2-ynyl ester	290 C18H26O3	1000299-24-1	35
3	3-Hydroxypyridine monoacetate	137 C7H7NO2	017747-43-2	35
4	Ethane-1,2-diamine, N,N'-bis(3-p...	248 C10H12N6O2	1000277-19-4	32
5	4-Pyridinol	95 C5H5NO	000626-64-2	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
Data File : BN004316.D
Acq On : 03 Jan 2019 13:58
Operator : JU/SJ
Sample : J6593-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampled :
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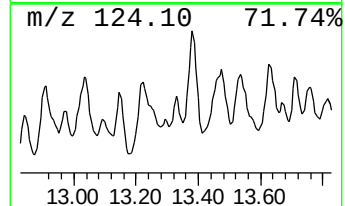
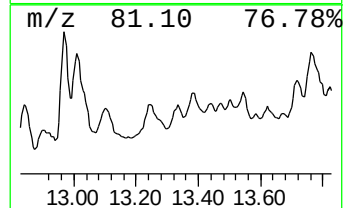
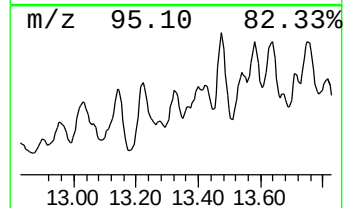
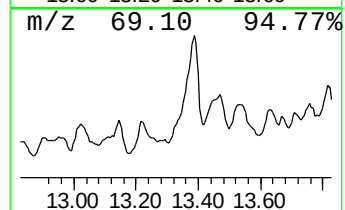
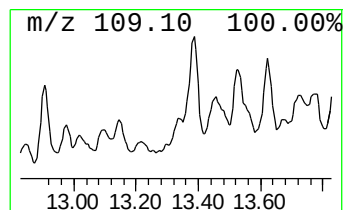
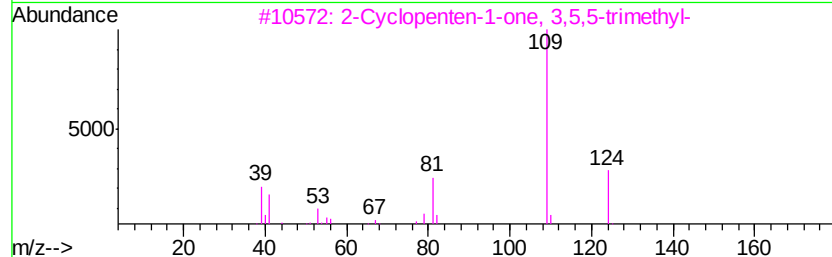
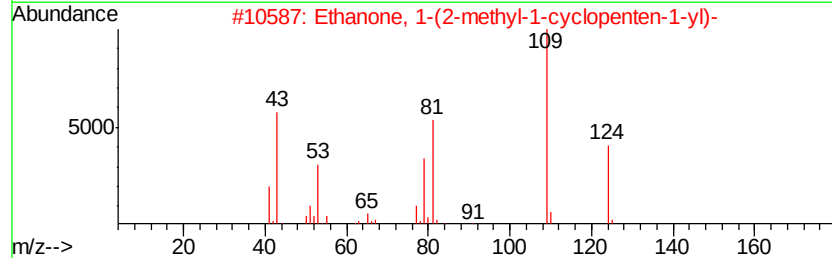
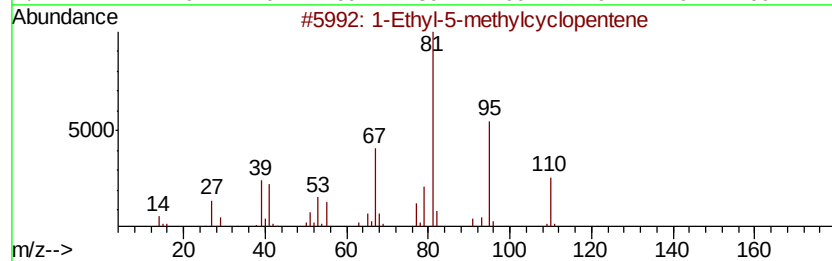
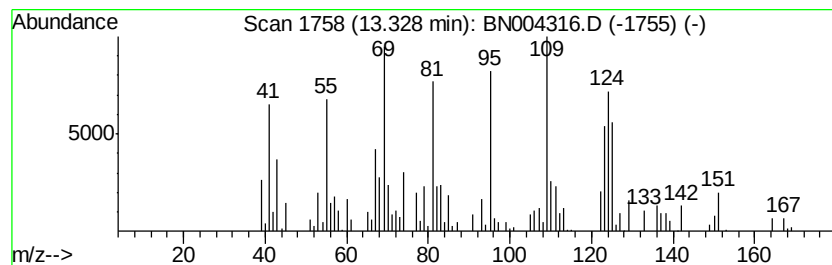
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 unknown-20 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	2.83 ng/ul	80298	Acenaphthene-d10	14.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	35
2		Ethanone, 1-(2-methyl-1-cyclopenten-1-yl)-	124	C8H12O	003168-90-9	27
3		2-Cyclopenten-1-one, 3,5,5-trimethyl-	124	C8H12O	024156-95-4	18
4		di-t-Butylacetylene	138	C10H18	017530-24-4	18
5		Methyl ethyl cyclopentene	110	C8H14	019780-56-4	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN010319\
 Data File : BN004316.D
 Acq On : 03 Jan 2019 13:58
 Operator : JU/SJ
 Sample : J6593-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.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
Indane	8.17	8.2	ng/ul	96187	1	7.81	234748	20.0
Triethyl phosphate	9.27	15.2	ng/ul	298975	2	10.60	392929	20.0
Pentanoic acid, 2...	10.16	6.5	ng/ul	126691	2	10.60	392929	20.0
unknown-01	10.55	6.5	ng/ul	128289	2	10.60	392929	20.0
unknown-02	10.79	4.5	ng/ul	87500	2	10.60	392929	20.0
unknown-03	10.90	5.7	ng/ul	111471	2	10.60	392929	20.0
unknown-04	10.96	9.8	ng/ul	193144	2	10.60	392929	20.0
unknown-05	11.10	21.6	ng/ul	424088	2	10.60	392929	20.0
unknown-06	11.19	4.7	ng/ul	91521	2	10.60	392929	20.0
unknown-07	11.24	5.6	ng/ul	109327	2	10.60	392929	20.0
Cyclopentene, 1,2...	11.40	10.5	ng/ul	206371	2	10.60	392929	20.0
Thiophane, propyl-	11.55	26.6	ng/ul	521745	2	10.60	392929	20.0
unknown-08	11.62	4.5	ng/ul	89455	2	10.60	392929	20.0
1-Butene, 1-(meth...	11.67	12.6	ng/ul	246543	2	10.60	392929	20.0
2-Propanol, 1-[2-...	11.78	14.1	ng/ul	276130	2	10.60	392929	20.0
unknown-09	12.03	6.0	ng/ul	117081	2	10.60	392929	20.0
Methane, diethoxy-	12.06	20.0	ng/ul	393252	2	10.60	392929	20.0
2-Propanol, 1-[1-...	12.12	27.6	ng/ul	542756	2	10.60	392929	20.0
unknown-10	12.30	20.7	ng/ul	405982	2	10.60	392929	20.0
unknown-11	12.38	3.7	ng/ul	72265	2	10.60	392929	20.0
unknown-12	12.45	8.2	ng/ul	161856	2	10.60	392929	20.0
unknown-13	12.61	14.5	ng/ul	411960	3	14.45	568396	20.0
unknown-14	12.71	9.4	ng/ul	268267	3	14.45	568396	20.0
unknown-15	12.93	17.6	ng/ul	500544	3	14.45	568396	20.0
unknown-16	12.97	3.7	ng/ul	105912	3	14.45	568396	20.0
unknown-17	13.02	6.8	ng/ul	193957	3	14.45	568396	20.0
unknown-18	13.14	3.5	ng/ul	100426	3	14.45	568396	20.0
unknown-19	13.22	6.5	ng/ul	184565	3	14.45	568396	20.0
unknown-20	13.33	2.8	ng/ul	80298	3	14.45	568396	20.0
unknown-21	13.39	10.6	ng/ul	302205	3	14.45	568396	20.0