

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
 Data File : BN012594.D
 Acq On : 18 Nov 2020 14:45
 Operator : CG/JU
 Sample : L4726-11
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBGE3

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 0.1 % 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-BN102220.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	2.523	3	6	24	rVB3	75672817	128132908	100.00%	44.973%
2	2.723	36	40	45	rBV	578660	687298	0.54%	0.241%
3	2.764	45	47	57	rVB	113678	142162	0.11%	0.050%
4	4.611	357	361	372	rVB	88423	139495	0.11%	0.049%
5	6.634	698	705	708	rBV	849033	1397470	1.09%	0.490%
6	6.799	727	733	742	rBV	608681	931036	0.73%	0.327%
7	6.887	743	748	758	rVV	1090841	1642332	1.28%	0.576%
8	6.987	758	765	775	rVB	866357	1394793	1.09%	0.490%
9	7.446	835	843	851	rBV	627943	973505	0.76%	0.342%
10	7.899	913	920	931	rBV	1925379	2969733	2.32%	1.042%
11	8.163	958	965	970	rBV	976383	1579392	1.23%	0.554%
12	8.222	970	975	979	rVV	2047629	3172488	2.48%	1.113%
13	8.263	979	982	1003	rVB	1584260	2436659	1.90%	0.855%
14	8.505	1017	1023	1033	rVB	960697	1543275	1.20%	0.542%
15	8.599	1033	1039	1048	rBV	797492	1295273	1.01%	0.455%
16	9.163	1126	1135	1150	rBV	1478312	2298927	1.79%	0.807%
17	9.310	1153	1160	1169	rBV	727877	1196210	0.93%	0.420%
18	9.416	1169	1178	1193	rVB	1572859	2485713	1.94%	0.872%
19	9.652	1211	1218	1231	rBV	1230463	1836396	1.43%	0.645%
20	9.846	1244	1251	1271	rBV2	1061792	2566504	2.00%	0.901%
21	10.216	1303	1314	1318	rBV	818939	1402159	1.09%	0.492%
22	10.263	1318	1322	1332	rVB	1052933	1659553	1.30%	0.582%
23	10.363	1332	1339	1353	rBV	522948	972855	0.76%	0.341%
24	10.546	1364	1370	1384	rVB	1992859	3187863	2.49%	1.119%
25	11.157	1466	1474	1495	rBV	531683	1514701	1.18%	0.532%
26	11.757	1569	1576	1583	rBV	796475	1252961	0.98%	0.440%
27	12.116	1629	1637	1652	rVB	2341631	3816101	2.98%	1.339%
28	12.269	1652	1663	1673	rVB2	1700916	4087712	3.19%	1.435%
29	12.587	1710	1717	1733	rBV	1452141	2344781	1.83%	0.823%
30	12.963	1773	1781	1793	rBV	1038868	1598153	1.25%	0.561%
31	13.528	1871	1877	1881	rBV	2032024	2698310	2.11%	0.947%
32	13.575	1881	1885	1893	rVB	1153764	1546446	1.21%	0.543%
33	13.793	1916	1922	1925	rBV	2089680	3221520	2.51%	1.131%
34	13.822	1925	1927	1936	rVB	1369461	1675457	1.31%	0.588%

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35	14.028	1955	1962	1969	rBV	1345683	2050674	1.60%	0.720%
36	14.104	1969	1975	1981	rVV	1299587	1898211	1.48%	0.666%
37	14.169	1981	1986	1992	rVV	2373538	3413882	2.66%	1.198%
38	14.234	1992	1997	2007	rVV	1184947	1720830	1.34%	0.604%
39	14.328	2007	2013	2022	rVV	907113	1461064	1.14%	0.513%
40	14.416	2022	2028	2038	rVV	1210145	1663457	1.30%	0.584%
41	14.510	2038	2044	2050	rVB	1245563	1733353	1.35%	0.608%
42	15.104	2139	2145	2150	rBV	2573299	3690951	2.88%	1.295%
43	15.163	2150	2155	2158	rVV	2745738	3756890	2.93%	1.319%
44	15.198	2158	2161	2165	rVV	1754322	2664107	2.08%	0.935%
45	15.257	2165	2171	2183	rVV3	3041585	5527816	4.31%	1.940%
46	16.510	2378	2384	2403	rBV	3862224	5599226	4.37%	1.965%
47	16.857	2437	2443	2454	rVB	1583245	2166691	1.69%	0.760%
48	16.957	2454	2460	2463	rBV	2952280	4214342	3.29%	1.479%
49	16.992	2463	2466	2480	rVB	3075813	3914965	3.06%	1.374%
50	17.269	2505	2513	2534	rBV	2978933	4314219	3.37%	1.514%
51	17.816	2602	2606	2616	rVB3	81857	148582	0.12%	0.052%
52	18.092	2649	2653	2659	rVB	145132	193950	0.15%	0.068%
53	18.928	2785	2795	2803	rBV	1966745	2674914	2.09%	0.939%
54	19.269	2847	2853	2862	rBV	3440449	4811979	3.76%	1.689%
55	20.222	3010	3015	3024	rBV	2531482	2897185	2.26%	1.017%
56	20.763	3102	3107	3112	rBV	153019	270490	0.21%	0.095%
57	20.816	3112	3116	3121	rVV	210746	312422	0.24%	0.110%
58	20.869	3121	3125	3133	rVV	614753	800991	0.63%	0.281%
59	21.010	3144	3149	3154	rBV	3378438	3529824	2.75%	1.239%
60	21.075	3155	3160	3163	rVV	1796116	2477637	1.93%	0.870%
61	21.110	3163	3166	3175	rVB	3973629	4721207	3.68%	1.657%
62	21.810	3281	3285	3291	rBV	698740	906335	0.71%	0.318%
63	22.104	3333	3335	3342	rVB3	119112	134227	0.10%	0.047%
64	22.610	3413	3421	3432	rVB	3692907	5903633	4.61%	2.072%
65	23.063	3492	3498	3502	rBV	2695056	4910873	3.83%	1.724%
66	23.110	3502	3506	3513	rVV	3505707	5674905	4.43%	1.992%
67	23.192	3515	3520	3531	rVB	1379608	2437671	1.90%	0.856%
68	25.280	3868	3875	3889	rVB	987890	2514274	1.96%	0.882%

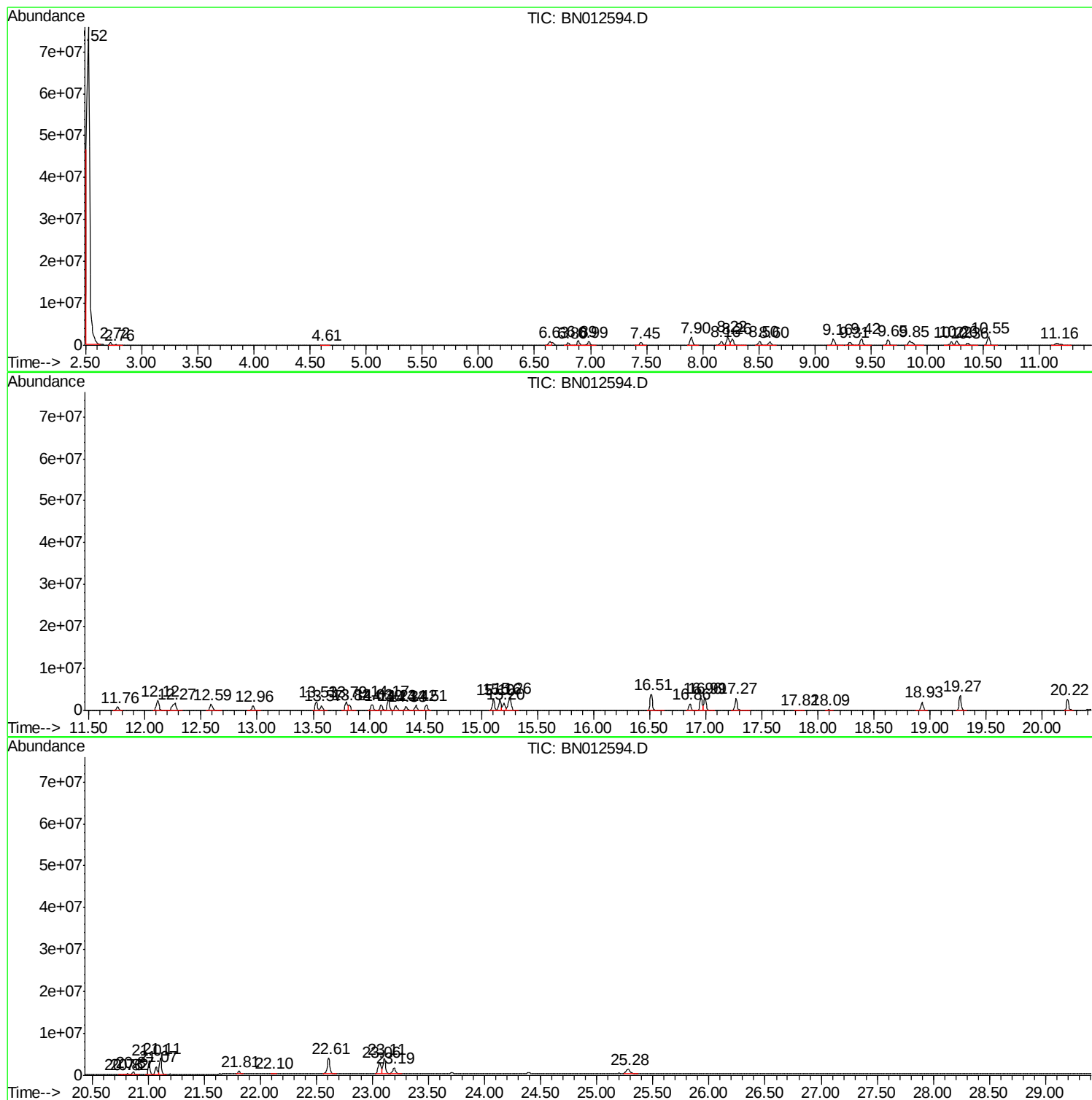
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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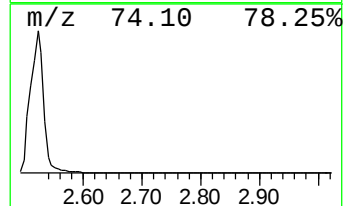
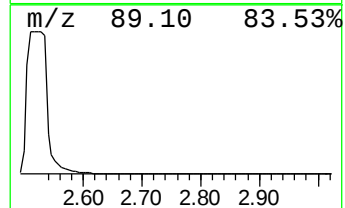
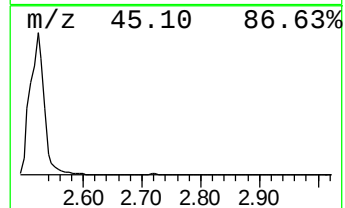
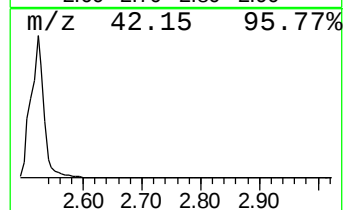
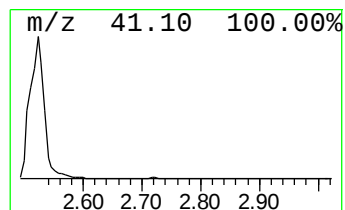
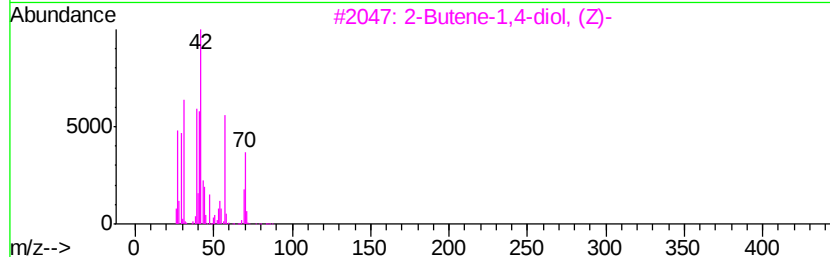
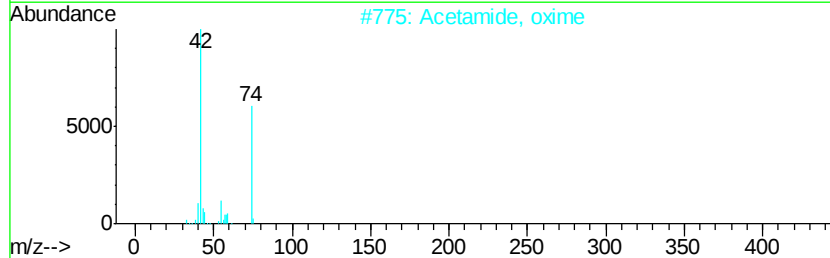
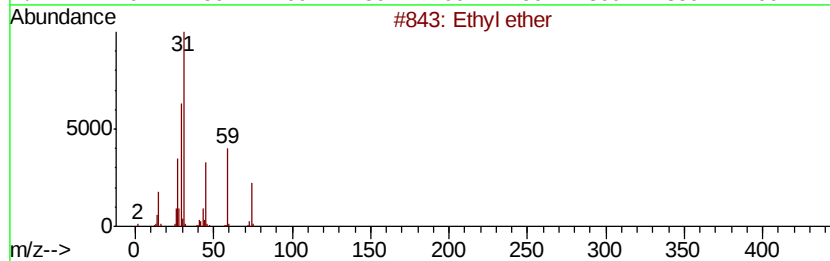
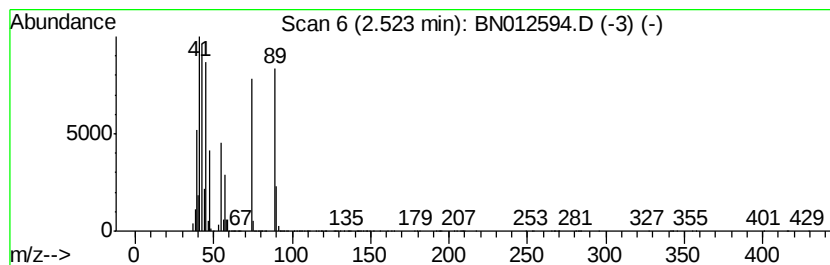
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN102220.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
2.52	2632.40 ng/ul	128133000	1,4-Dichlorobenzene-d4	7.45		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethyl ether	74	C4H10O	000060-29-7	9
2		Acetamide, oxime	74	C2H6N2O	022059-22-9	9
3		2-Butene-1,4-diol, (Z)-	88	C4H8O2	006117-80-2	9
4		2-Propanol, 1,1'-oxybis-	134	C6H14O3	000110-98-5	9
5		Propanoic acid	74	C3H6O2	000079-09-4	7



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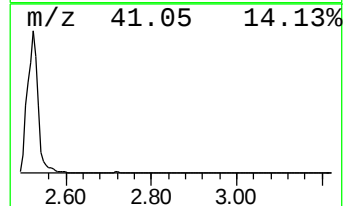
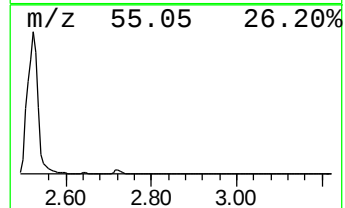
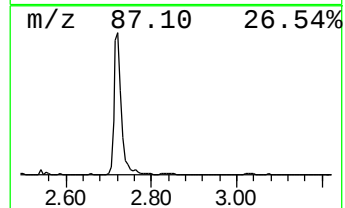
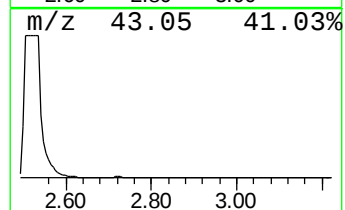
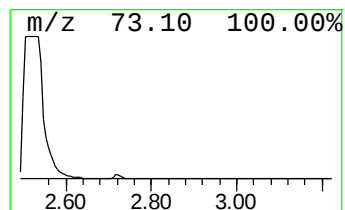
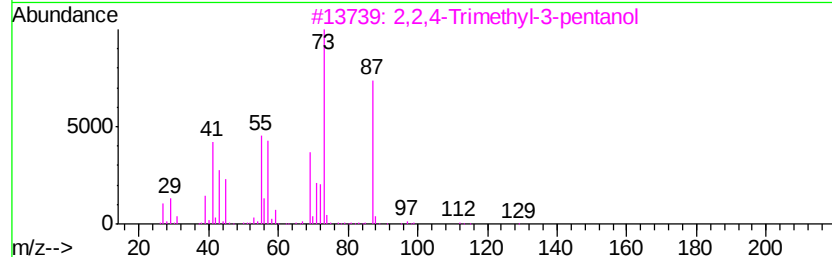
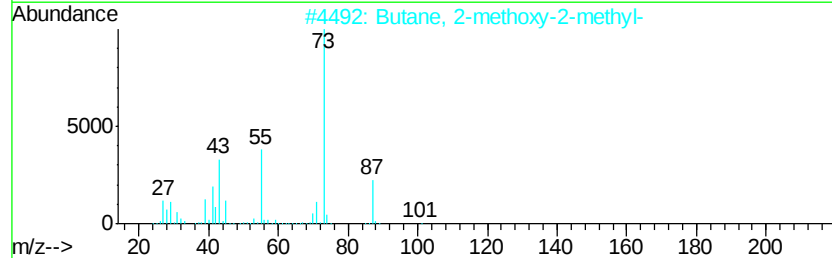
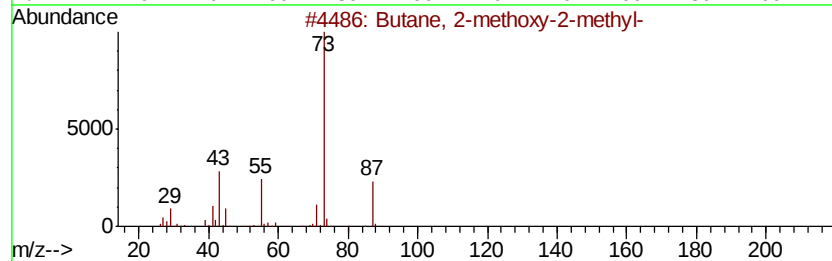
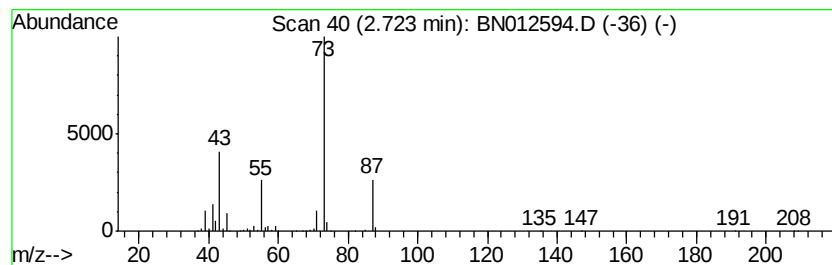
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.72	14.12 ng/ul	687298	1,4-Dichlorobenzene-d4	7.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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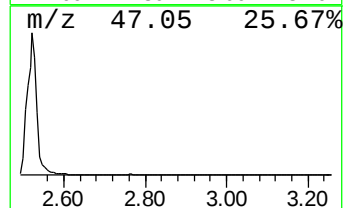
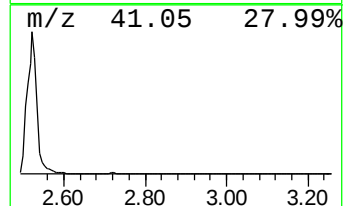
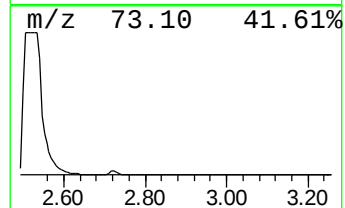
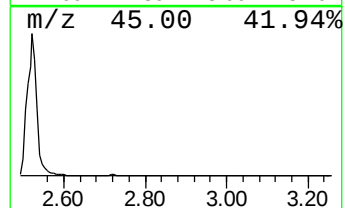
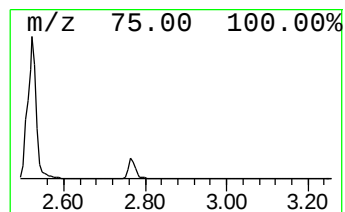
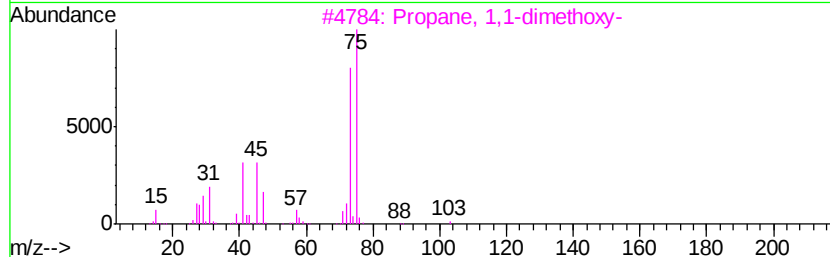
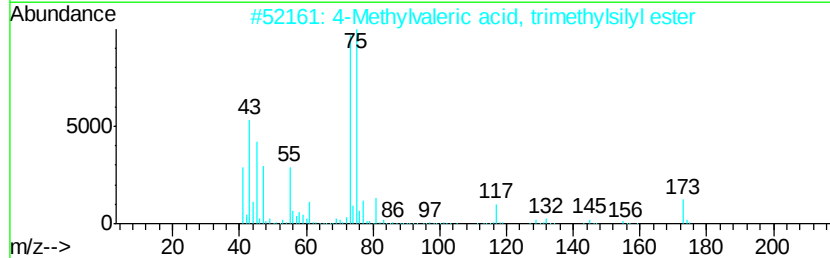
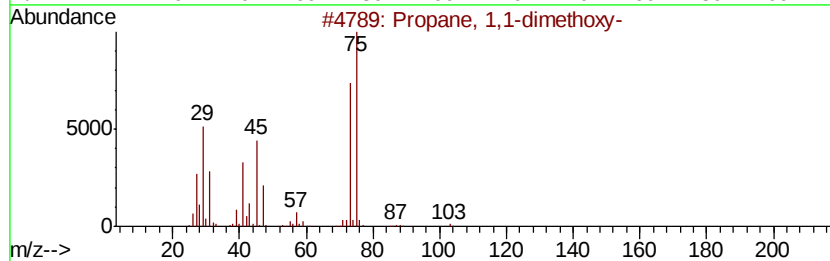
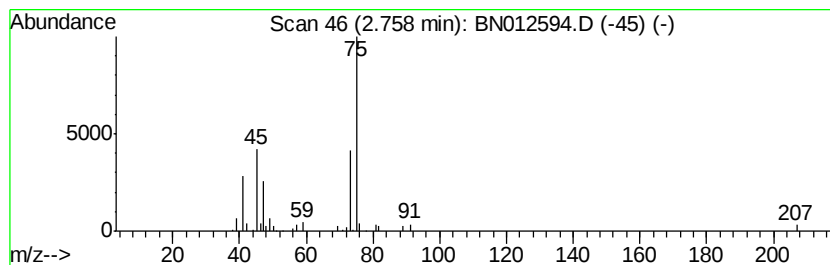
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Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.76	2.92 ng/ul	142162	1,4-Dichlorobenzene-d4	7.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	64
2		4-Methylvaleric acid, trimethyls...	188	C9H20O2Si	959048-10-3	40
3		Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	40
4		Acetic acid, dimethoxy-, methyl ...	134	C5H10O4	000089-91-8	40
5		3-Bromopropionaldehyde dimethyl ...	182	C5H11BrO2	036255-44-4	38



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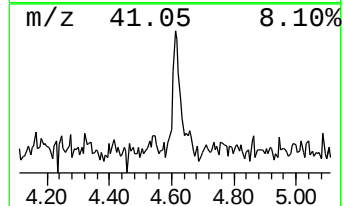
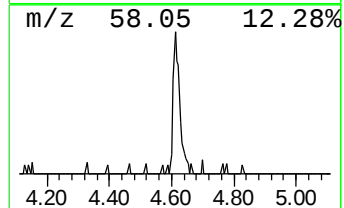
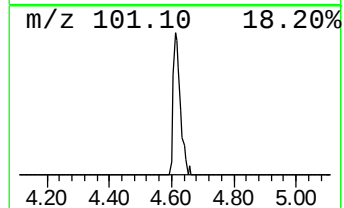
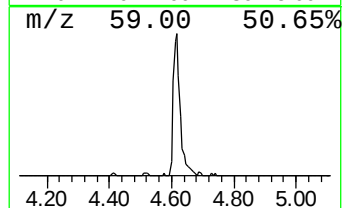
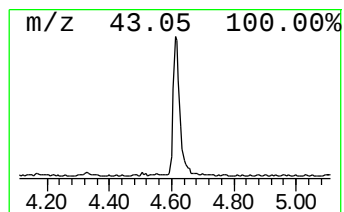
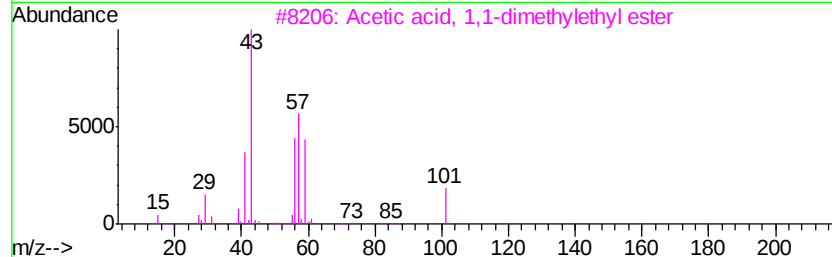
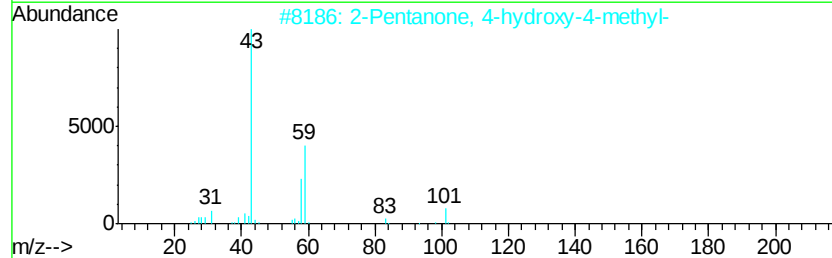
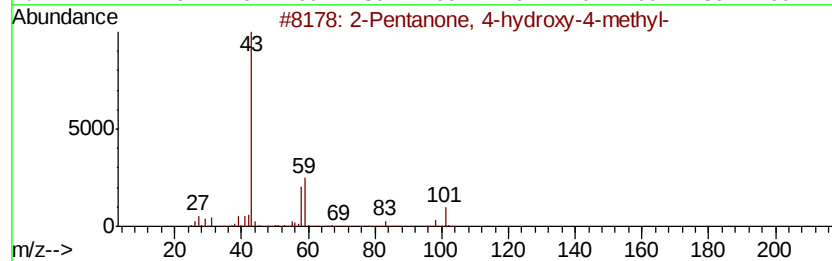
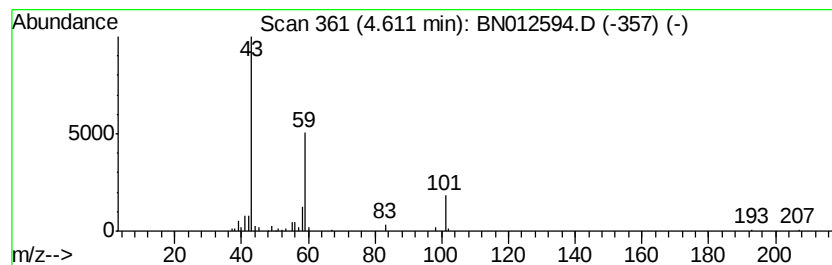
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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.61	2.87 ng/ul	139495	1,4-Dichlorobenzene-d4	7.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	43
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
Operator : CG/JU
Sample : L4726-11
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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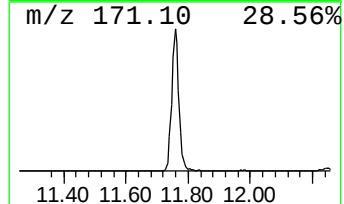
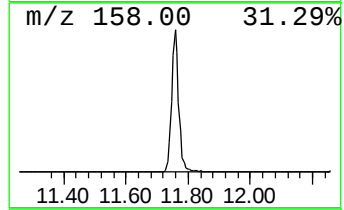
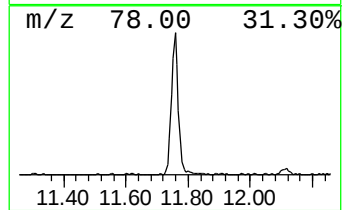
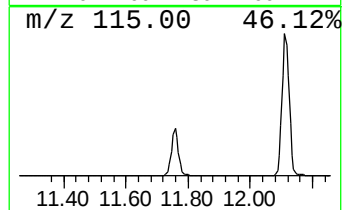
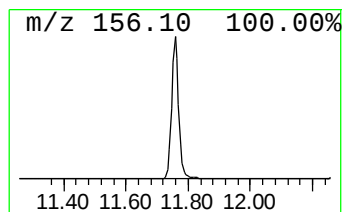
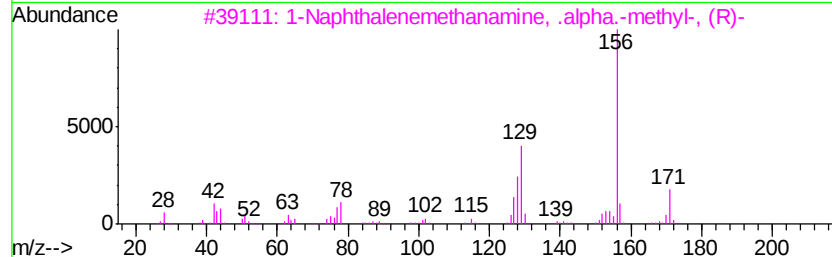
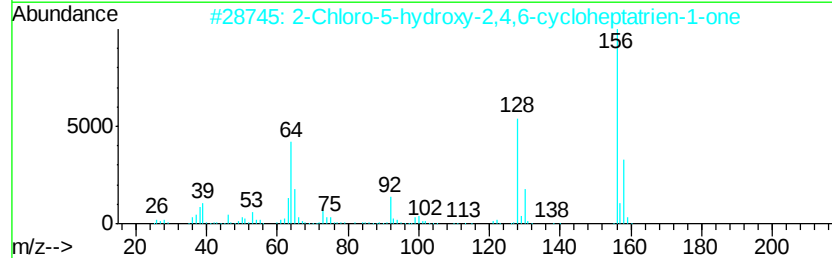
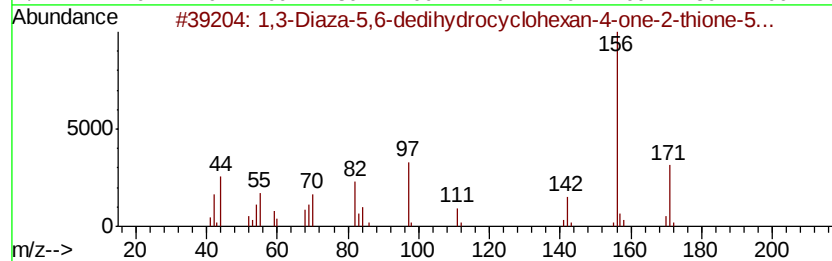
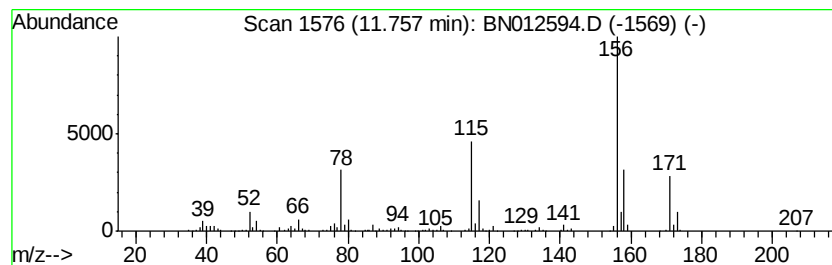
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	17.87 ng/ul	1252960	Naphthalene-d8	10.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Diaza-5,6-dedihydrocyclohexa...	171	C6H9N3OS	021263-85-4	38
2		2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	38
3		1-Naphthalenemethanamine, .alpha...	171	C12H13N	003886-70-2	38
4		Tebuthiuron	228	C9H16N4OS	034014-18-1	37
5		Tebuthiuron	228	C9H16N4OS	034014-18-1	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
Operator : CG/JU
Sample : L4726-11
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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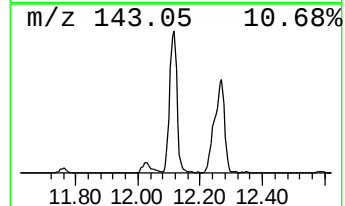
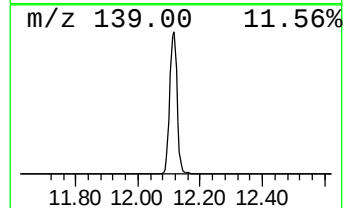
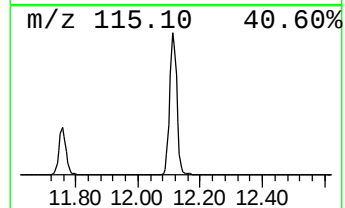
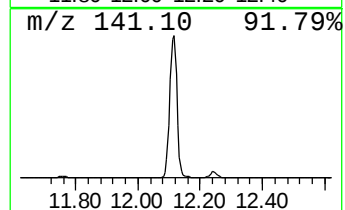
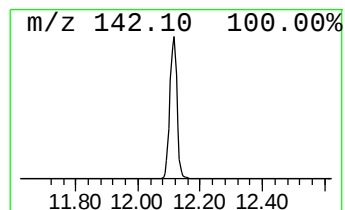
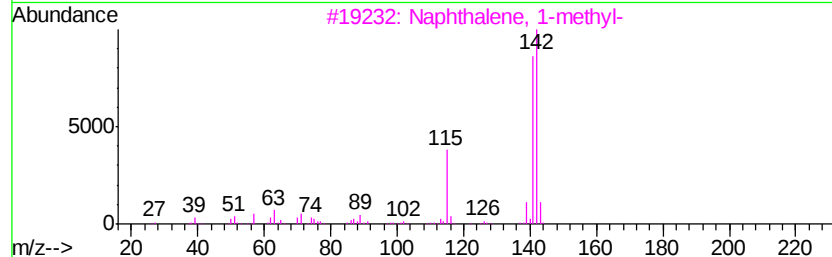
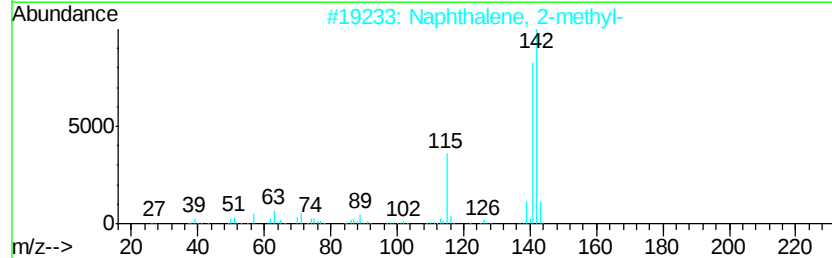
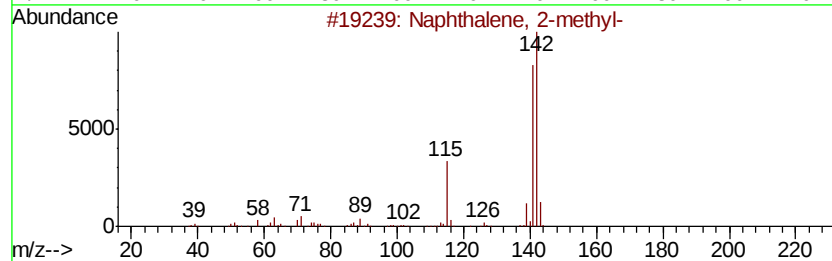
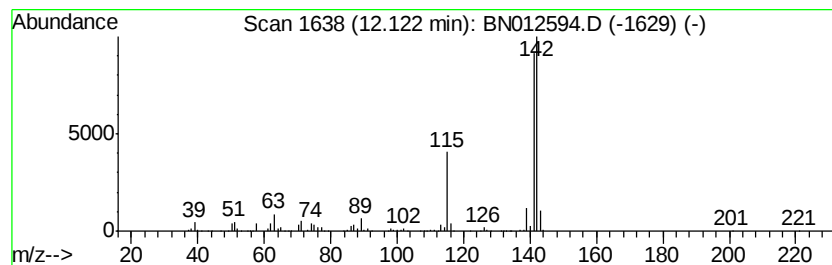
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	54.43 ng/ul	3816100	Naphthalene-d8	10.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
Operator : CG/JU
Sample : L4726-11
Misc :
ALS Vial : 4 Sample Multiplier: 1

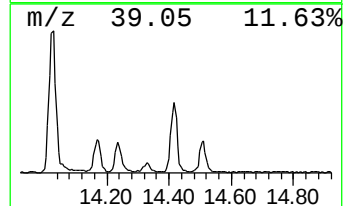
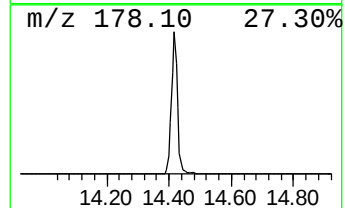
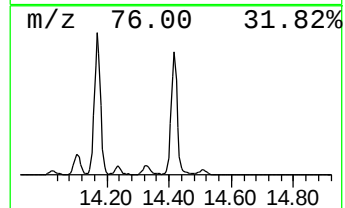
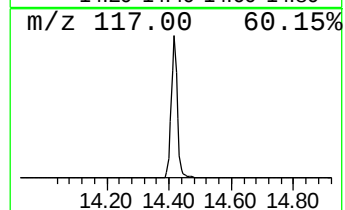
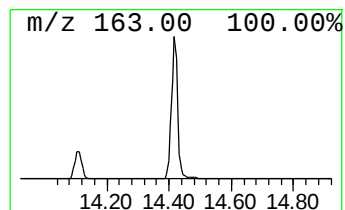
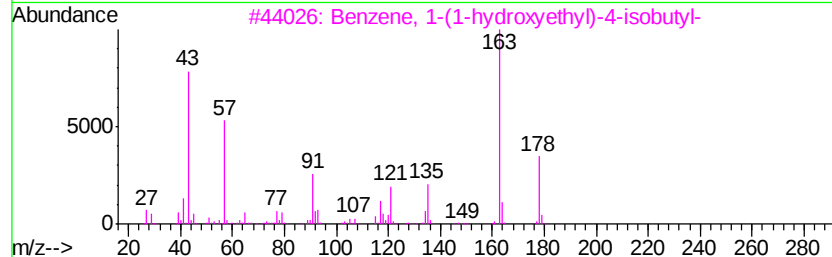
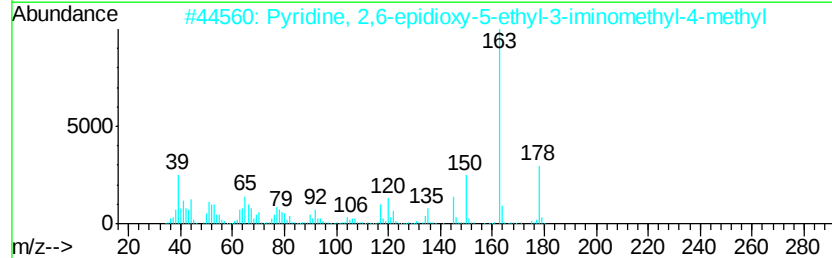
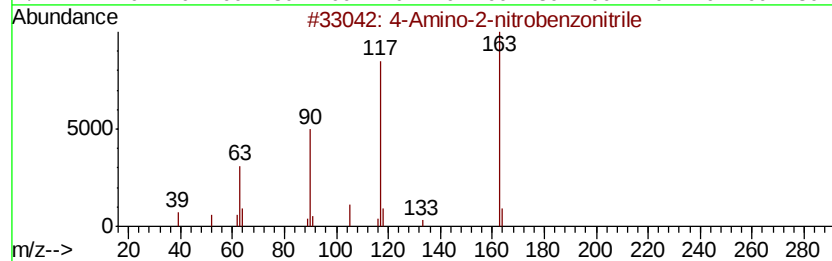
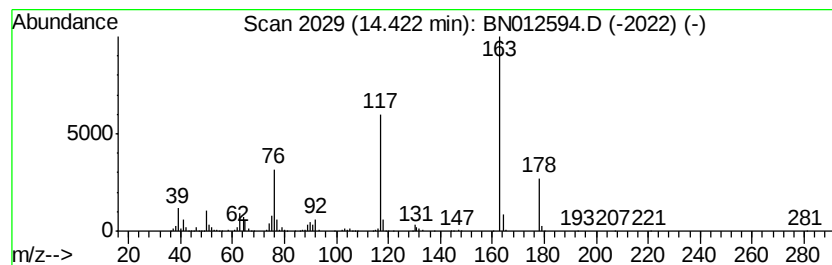
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN102220.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 4-Amino-2-nitrobenzonitrile Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.42	17.53 ng/ul	1663460	Acenaphthene-d10	14.10		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Amino-2-nitrobenzonitrile	163	C7H5N3O2	072115-07-2	53	
2	Pyridine, 2,6-epidioxv-5-ethyl-3...	178	C9H10N2O2	1000263-31-1	50	
3	Benzene, 1-(1-hvdroxvethyl)-4-is...	178	C12H18O	040150-92-3	47	
4	Propofol	178	C12H18O	002078-54-8	47	
5	Methyl p-chlorophenylethynyl ketone	178	C10H7ClO	039180-26-2	47	



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
Operator : CG/JU
Sample : L4726-11
Misc :
ALS Vial : 4 Sample Multiplier: 1

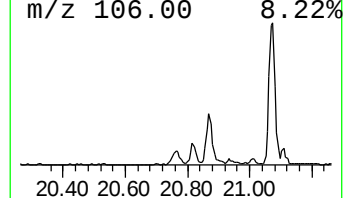
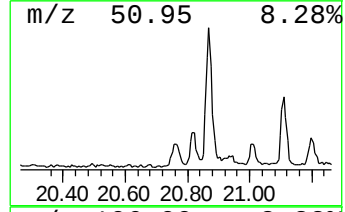
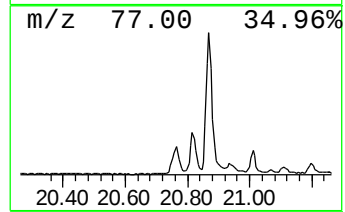
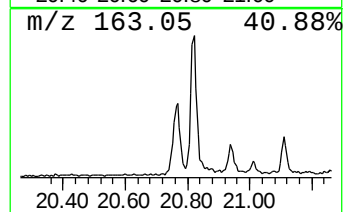
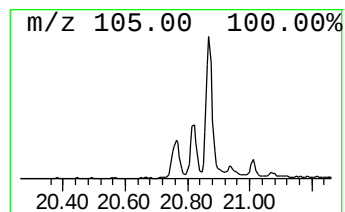
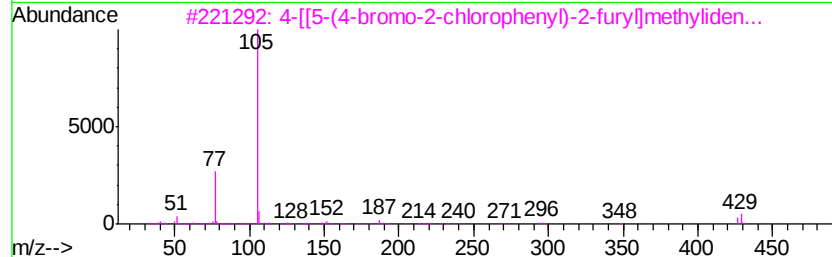
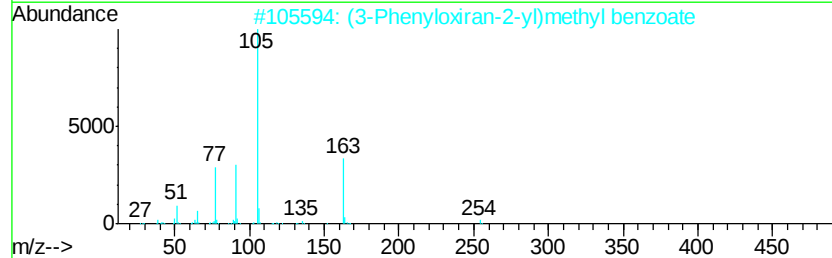
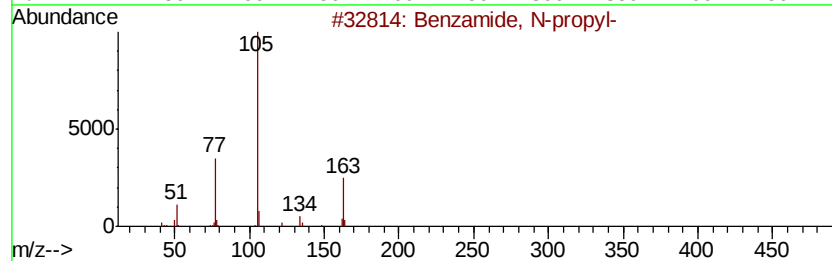
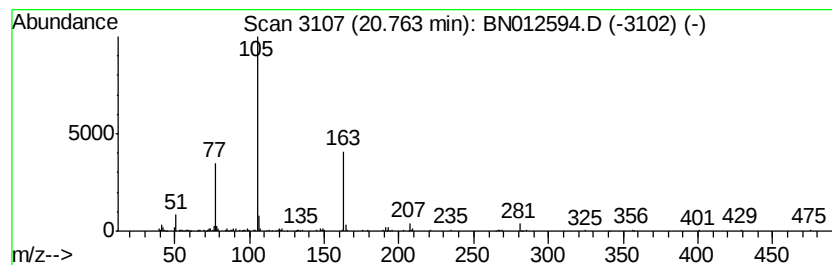
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzamide, N-propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.76	2.18 ng/ul	270490	Chrysene-d12	21.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzamide, N-propyl-	163 C10H13NO	010546-70-0 64
2		(3-Phenyloxiran-2-yl)methyl benz...	254 C16H14O3	1000366-96-1 50
3		4-[[5-(4-bromo-2-chlorophenyl)-2...	427 C20H11BrClNO3	282539-13-3 35
4		4-(((5-Methyl-3-isoxazolyl)amino...	269 C14H11N3O3	1000332-11-5 35
5		2,3,4-O-Tribenzoyl-1-ribose...	464 C26H21F07	1000129-81-2 27



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
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Sample : L4726-11
Misc :
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Instrument :
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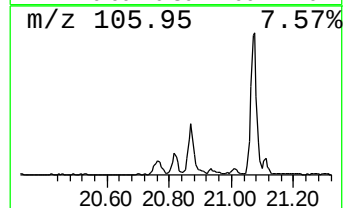
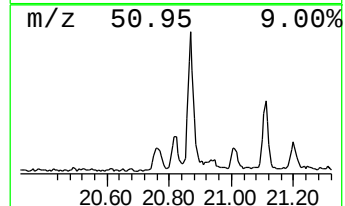
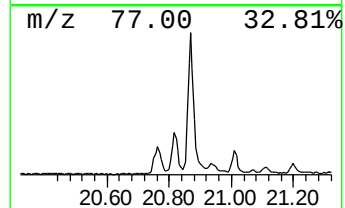
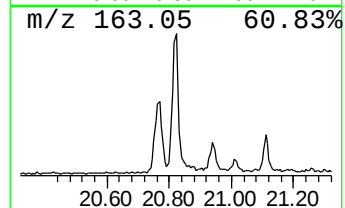
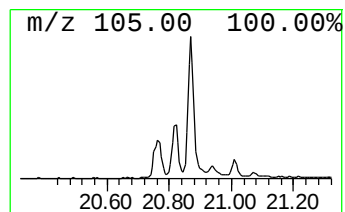
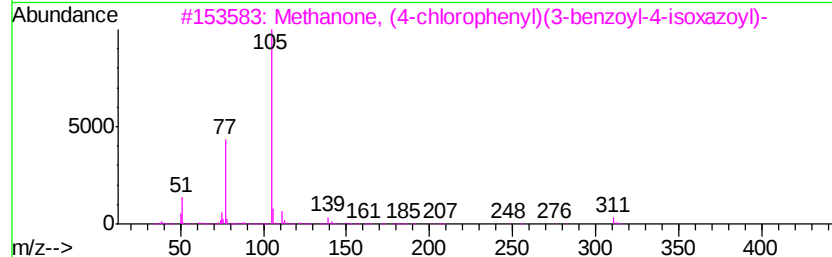
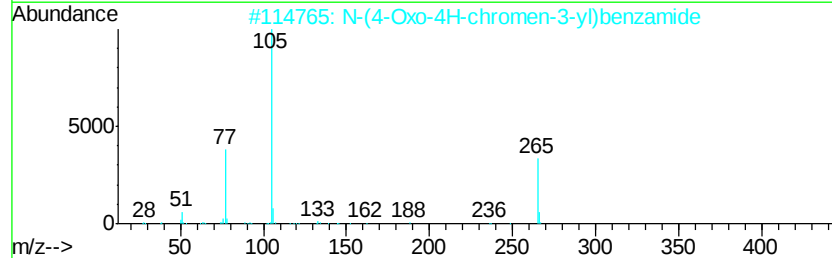
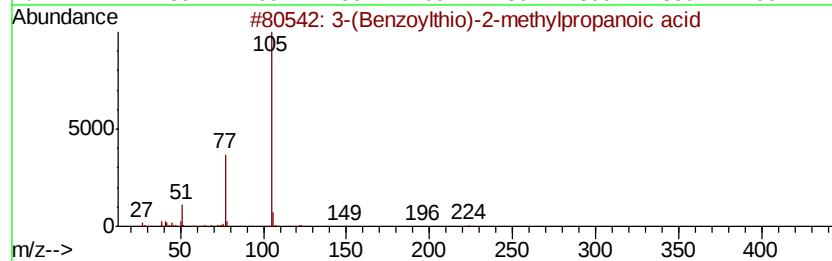
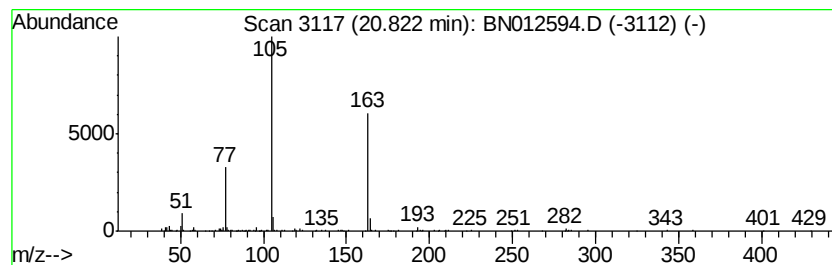
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 3-(Benzoylthio)-2-methylpro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	2.52 ng/ul	312422	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-(Benzoylthio)-2-methylpropanoi...	224	C11H12O3S	067714-34-5	60
2		N-(4-Oxo-4H-chromen-3-yl)benzamide	265	C16H11N03	1000243-49-5	49
3		Methanone, (4-chlorophenyl)(3-be...	311	C17H10ClN03	1000276-36-3	47
4		Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	47
5		Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
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Misc :
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ClientSampleId :
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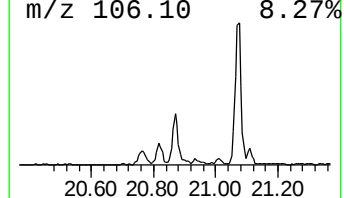
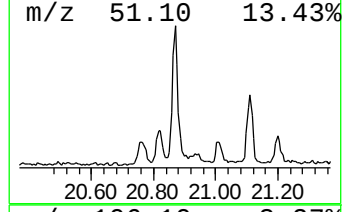
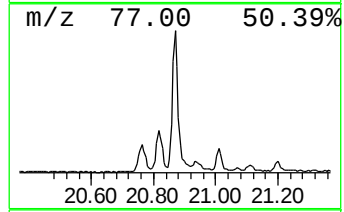
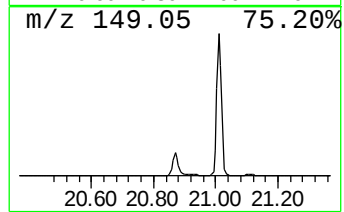
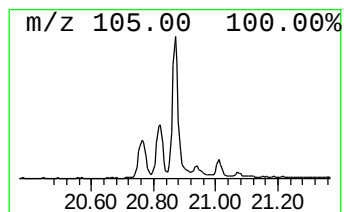
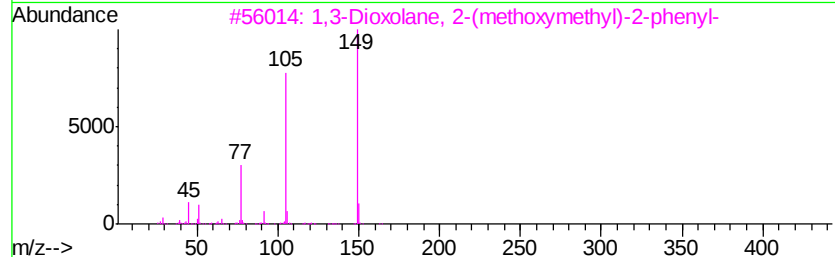
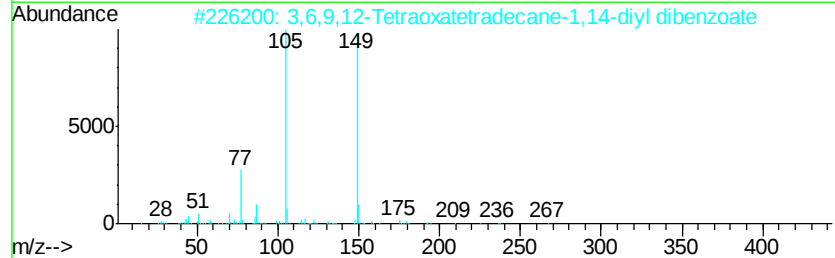
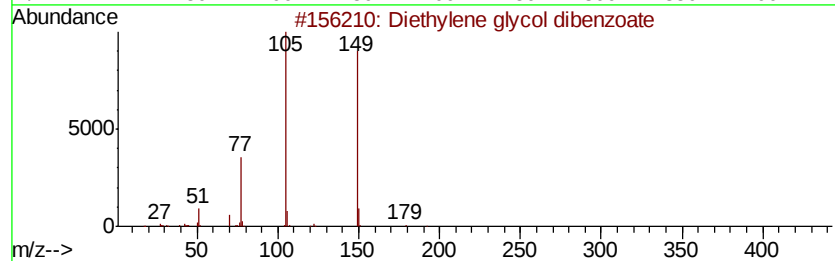
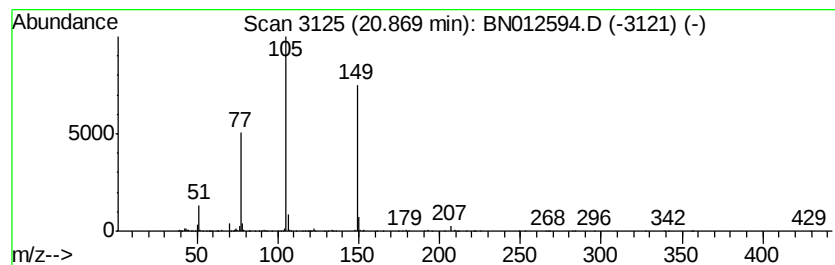
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Diethylene glycol dibenzoate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	6.47 ng/ul	800991	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	64
2		3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	64
3		1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	64
4		Vinyl benzoate	148	C9H8O2	000769-78-8	38
5		1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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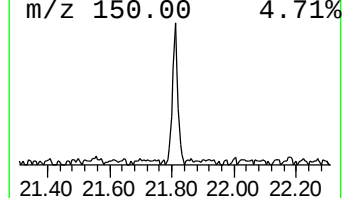
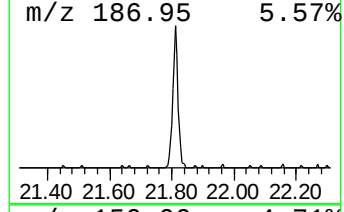
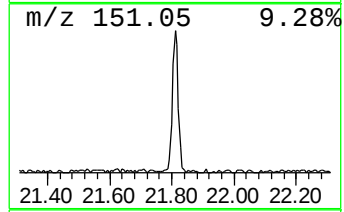
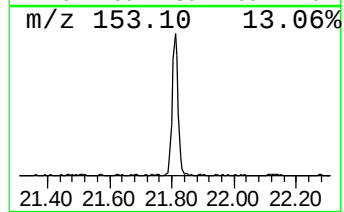
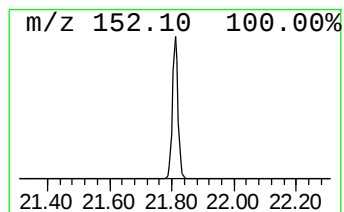
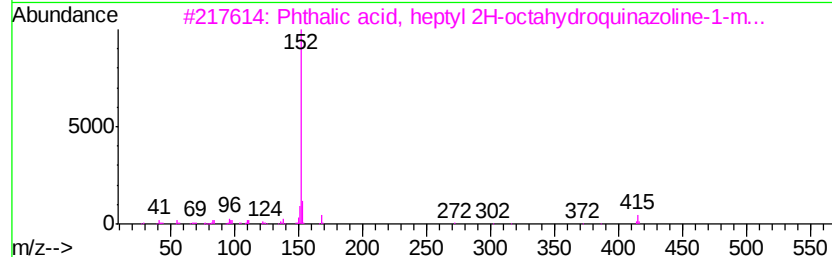
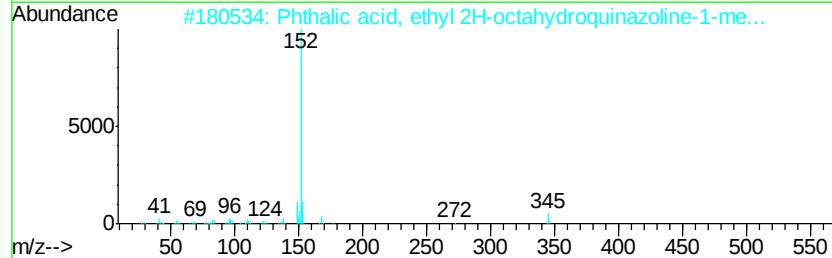
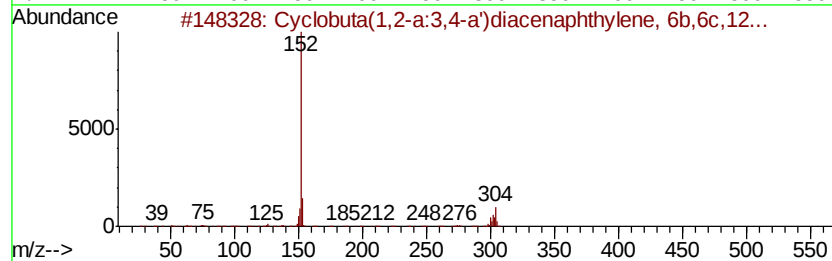
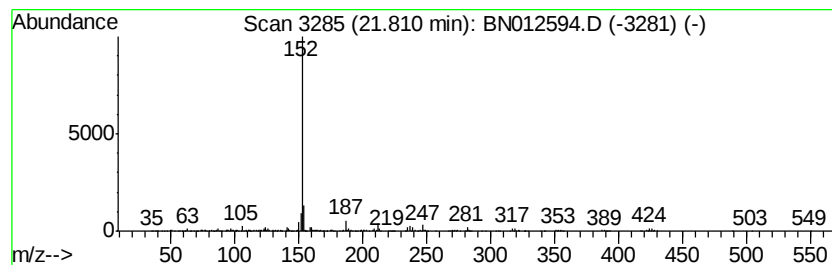
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN102220.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 Cyclobuta(1,2-a:3,4-a')diac... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	7.32 ng/ul	906335	Chrysene-d12	21.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclobuta(1,2-a:3,4-a')diacenaph...	304	C24H16	007259-03-2	80
2		Phthalic acid, ethyl 2H-octahydr...	345	C20H27N04	1000322-80-1	56
3		Phthalic acid, heptyl 2H-octahyd...	415	C25H37N04	1000322-80-8	56
4		Acenaphthylene	152	C12H8	000208-96-8	45
5		Biphenylene	152	C12H8	000259-79-0	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN111820\
Data File : BN012594.D
Acq On : 18 Nov 2020 14:45
Operator : CG/JU
Sample : L4726-11
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBGE3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN102220.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
unknown-01	2.52	2632.4	ng/ul	128133000	1	7.45	973505	20.0
Butane, 2-methoxy...	2.72	14.1	ng/ul	687298	1	7.45	973505	20.0
Propane, 1,1-dime...	2.76	2.9	ng/ul	142162	1	7.45	973505	20.0
2-Pentanone, 4-hy...	4.61	2.9	ng/ul	139495	1	7.45	973505	20.0
unknown-02	11.76	17.9	ng/ul	1252960	2	10.22	1402160	20.0
Naphthalene, 1-me...	12.12	54.4	ng/ul	3816100	2	10.22	1402160	20.0
4-Amino-2-nitrobe...	14.42	17.5	ng/ul	1663460	3	14.10	1898210	20.0
Benzamide, N-propyl-	20.76	2.2	ng/ul	270490	5	21.07	2477640	20.0
3-(Benzoylthio)-2...	20.82	2.5	ng/ul	312422	5	21.07	2477640	20.0
Diethylene glycol...	20.87	6.5	ng/ul	800991	5	21.07	2477640	20.0
Cyclobuta(1,2-a:3...	21.81	7.3	ng/ul	906335	5	21.07	2477640	20.0