

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
 Data File : BN008555.D
 Acq On : 14 Nov 2019 13:53
 Operator : JU
 Sample : K5656-10
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5458

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-BN111319MA.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	3.017	3	5	14	rVB	237469	272324	2.65%	0.100%
2	3.217	35	39	45	rVV	93654	117099	1.14%	0.043%
3	3.275	46	49	53	rVB	28957	34347	0.33%	0.013%
4	3.552	92	96	100	rBV2	40484	52903	0.52%	0.020%
5	3.846	141	146	149	rVB	19429	30130	0.29%	0.011%
6	3.922	156	159	168	rVB	24863	38334	0.37%	0.014%
7	4.240	208	213	218	rBV3	22039	33128	0.32%	0.012%
8	4.440	241	247	255	rVB10	17720	37012	0.36%	0.014%
9	4.811	306	310	317	rBV	278002	399856	3.90%	0.148%
10	5.864	484	489	496	rVB	37861	67125	0.65%	0.025%
11	6.411	577	582	586	rVB2	21782	30742	0.30%	0.011%
12	6.816	643	651	654	rBV	1719723	3085547	30.07%	1.139%
13	6.975	671	678	687	rBV	1189997	1851034	18.04%	0.683%
14	7.063	687	693	704	rVB	2246031	3316006	32.32%	1.224%
15	7.175	704	712	719	rBV	1637897	2681621	26.13%	0.989%
16	7.628	783	789	799	rBV	1320564	2062601	20.10%	0.761%
17	8.075	856	865	873	rBV	3756673	5706281	55.61%	2.106%
18	8.340	903	910	915	rBV	1977179	3212587	31.31%	1.185%
19	8.399	915	920	923	rVV	3822288	5776138	56.29%	2.131%
20	8.440	923	927	943	rVB	2631948	4247172	41.39%	1.567%
21	8.699	965	971	977	rBV	1708504	2832840	27.61%	1.045%
22	8.769	977	983	997	rVB	1581692	2566688	25.01%	0.947%
23	9.340	1072	1080	1088	rBV	2527882	4078999	39.75%	1.505%
24	9.487	1096	1105	1115	rBV	1296474	2148044	20.93%	0.793%
25	9.593	1115	1123	1130	rBV	2910056	4654609	45.36%	1.718%
26	9.822	1154	1162	1170	rVB	2360391	3534777	34.45%	1.304%
27	10.022	1189	1196	1211	rBV2	2075615	4967869	48.41%	1.833%
28	10.246	1228	1234	1243	rBV4	21130	42564	0.41%	0.016%
29	10.399	1251	1260	1264	rBV	1793202	3063507	29.86%	1.130%
30	10.446	1264	1268	1275	rVV	1956388	3101154	30.22%	1.144%
31	10.528	1275	1282	1292	rVV	144873	277883	2.71%	0.103%
32	10.634	1294	1300	1306	rVV7	20219	56739	0.55%	0.021%
33	10.693	1306	1310	1313	rVV	63768	98888	0.96%	0.036%
34	10.746	1313	1319	1330	rVB	3624635	5922057	57.71%	2.185%

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Integration Parameters: LSCINT.P

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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
 Title : SVOA CALIBRATION

35	11.116	1377	1382	1388	rBV	84471	147905	1.44%	0.055%
36	11.181	1388	1393	1400	rVB	151154	235978	2.30%	0.087%
37	11.304	1406	1414	1437	rBV	982109	2655227	25.88%	0.980%
38	11.528	1447	1452	1458	rVB4	20309	33492	0.33%	0.012%
39	11.934	1512	1521	1528	rBV	3486819	5644967	55.01%	2.083%
40	11.981	1528	1529	1536	rVB3	48736	68368	0.67%	0.025%
41	12.093	1538	1548	1554	rBV2	65298	120480	1.17%	0.044%
42	12.198	1560	1566	1573	rBV	324201	539473	5.26%	0.199%
43	12.287	1573	1581	1588	rVB	4431110	6871410	66.97%	2.536%
44	12.440	1598	1607	1614	rBV2	4068030	7082219	69.02%	2.613%
45	12.757	1651	1661	1669	rBV	2717239	4273835	41.65%	1.577%
46	13.122	1714	1723	1731	rVB	2008290	2919686	28.45%	1.077%
47	13.675	1811	1817	1821	rBV	3264773	4564746	44.49%	1.684%
48	13.722	1821	1825	1831	rVV	1912216	2596852	25.31%	0.958%
49	13.951	1855	1864	1866	rBV	3548820	5661251	55.17%	2.089%
50	13.975	1866	1868	1876	rVB	2376797	3102513	30.24%	1.145%
51	14.157	1889	1899	1906	rBV	1778749	2587672	25.22%	0.955%
52	14.257	1909	1916	1922	rVV	2538539	3867049	37.69%	1.427%
53	14.322	1922	1927	1931	rVV	4080760	5989263	58.37%	2.210%
54	14.363	1931	1934	1944	rVV	1553499	2101185	20.48%	0.775%
55	14.463	1944	1951	1961	rVV	1826173	2632135	25.65%	0.971%
56	14.563	1961	1968	1978	rVV	3031515	4131971	40.27%	1.525%
57	14.663	1978	1985	1993	rVB	2164903	3173029	30.92%	1.171%
58	15.045	2044	2050	2055	rBV7	36425	51974	0.51%	0.019%
59	15.128	2055	2064	2073	rBV	900770	1255089	12.23%	0.463%
60	15.257	2079	2086	2091	rBV	4533228	6418425	62.55%	2.368%
61	15.316	2091	2096	2102	rVV2	6196922	10261145	100.00%	3.786%
62	15.387	2102	2108	2121	rVV2	3902845	6755541	65.84%	2.493%
63	15.498	2121	2127	2131	rVB3	57106	90893	0.89%	0.034%
64	16.039	2216	2219	2222	rBV4	20054	22071	0.22%	0.008%
65	16.228	2246	2251	2260	rVB7	19674	33515	0.33%	0.012%
66	16.657	2318	2324	2336	rBV	5626863	8203342	79.95%	3.027%
67	16.910	2362	2367	2374	rBV	270574	371597	3.62%	0.137%
68	17.004	2377	2383	2393	rVB	3210755	4461168	43.48%	1.646%
69	17.104	2393	2400	2403	rBV	4841639	7305727	71.20%	2.696%
70	17.139	2403	2406	2412	rVB	4471446	5662429	55.18%	2.089%
71	17.269	2424	2428	2432	rVB6	27813	40759	0.40%	0.015%

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72	17.404	2444	2451	2458	rBV	5291642	7003063	68.25%	2.584%
73	17.510	2462	2469	2476	rBV	410713	576199	5.62%	0.213%
74	17.581	2478	2481	2485	rVV4	65117	98820	0.96%	0.036%
75	17.634	2485	2490	2498	rVB	459987	591702	5.77%	0.218%
76	17.828	2520	2523	2528	rVB5	24684	29814	0.29%	0.011%
77	17.963	2541	2546	2550	rVB	192208	278888	2.72%	0.103%
78	18.075	2561	2565	2569	rBV	68271	93193	0.91%	0.034%
79	18.181	2578	2583	2587	rBV	176757	243173	2.37%	0.090%
80	18.245	2588	2594	2599	rVV	271159	397834	3.88%	0.147%
81	18.298	2599	2603	2610	rVB9	33091	59838	0.58%	0.022%
82	18.375	2612	2616	2618	rBV5	38967	51212	0.50%	0.019%
83	18.416	2618	2623	2627	rVB	254755	331610	3.23%	0.122%
84	18.533	2641	2643	2647	rVB5	24872	27435	0.27%	0.010%
85	18.675	2663	2667	2671	rBV4	32271	48260	0.47%	0.018%
86	19.063	2723	2733	2740	rBV	3154869	4596203	44.79%	1.696%
87	19.286	2767	2771	2777	rBV	52373	85264	0.83%	0.031%
88	19.404	2784	2791	2798	rBV	5439799	7798514	76.00%	2.878%
89	20.357	2948	2953	2958	rBV	3869699	4346173	42.36%	1.604%
90	20.998	3059	3062	3066	rBV	209607	235686	2.30%	0.087%
91	21.157	3084	3089	3093	rBV	4806321	5277287	51.43%	1.947%
92	21.204	3093	3097	3100	rVV	3668140	4709365	45.90%	1.738%
93	21.239	3100	3103	3108	rVB	5854913	7110095	69.29%	2.624%
94	21.963	3222	3226	3232	rVB	951023	1271800	12.39%	0.469%
95	22.786	3360	3366	3377	rVB	5425068	8952420	87.25%	3.303%
96	23.251	3439	3445	3449	rBV	3765269	7219818	70.36%	2.664%
97	23.298	3449	3453	3458	rVV	4532794	7548321	73.56%	2.785%
98	23.345	3458	3461	3463	rVV	438221	630431	6.14%	0.233%
99	23.386	3463	3468	3477	rVB	2474142	4451411	43.38%	1.643%
100	25.545	3828	3835	3847	rVB	1821453	4610993	44.94%	1.701%

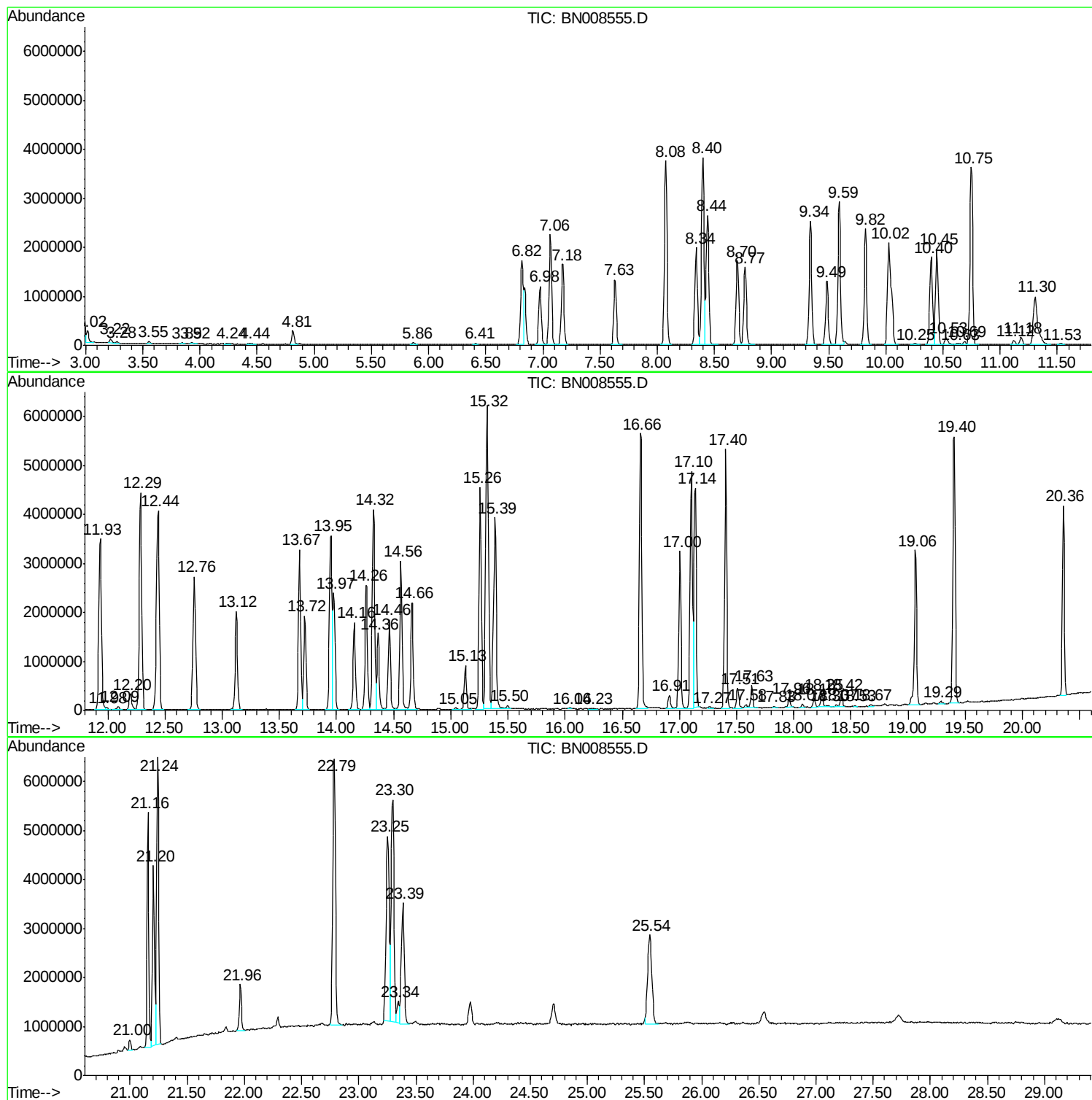
Sum of corrected areas: 271007808

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Instrument :
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

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



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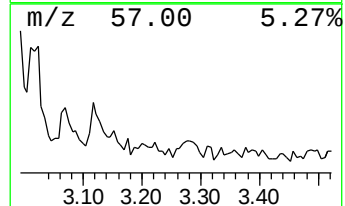
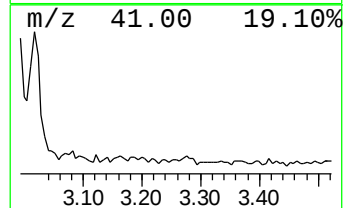
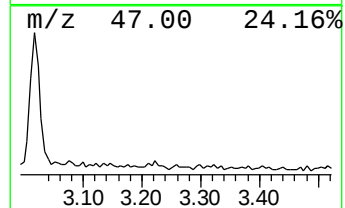
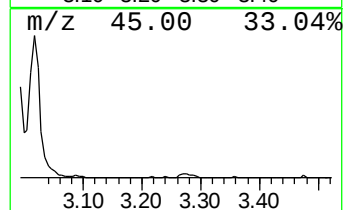
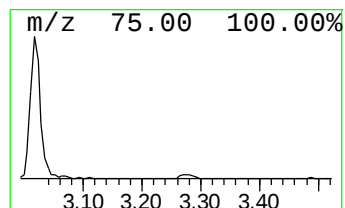
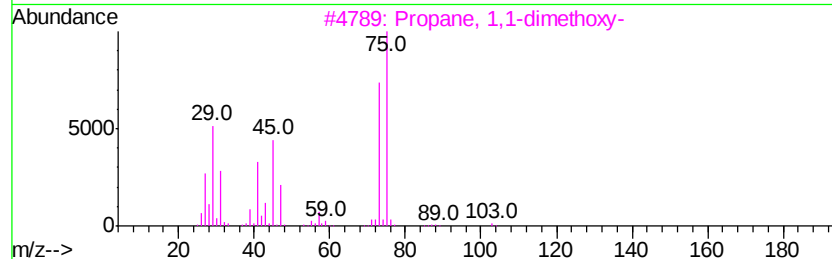
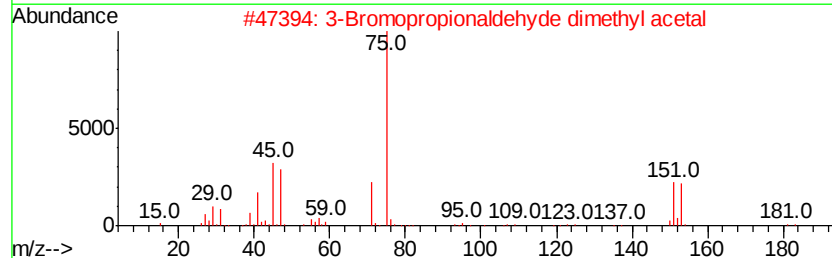
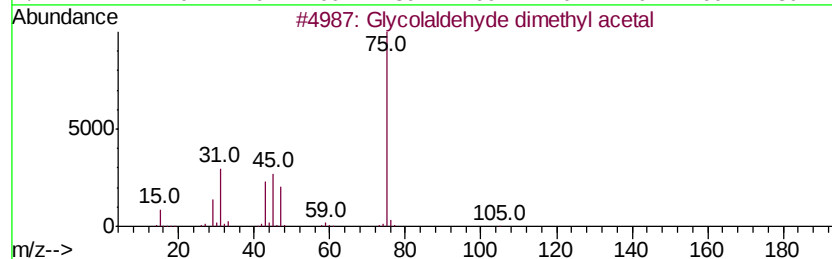
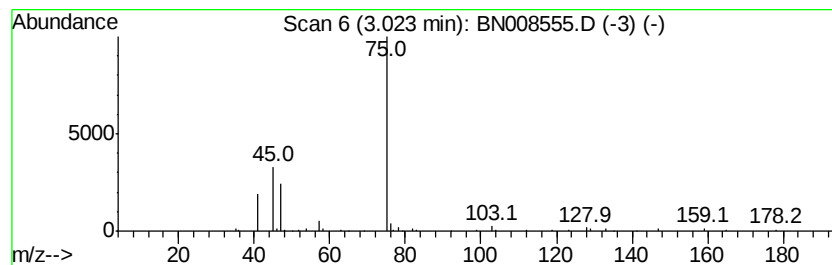
Instrument :
BNA_N
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A5458

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Glycolaldehyde dimethyl acetal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
3.02	2.64 ng/ul	272324	1,4-Dichlorobenzene-d4	7.63		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	50
2		3-Bromopropionaldehyde dimethyl ...	182	C5H11BrO2	036255-44-4	39
3		Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	39
4		Butane, 2-(methylthio)-	104	C5H12S	010359-64-5	9
5		Ethane, 1,1,2-trimethoxy-	120	C5H12O3	024332-20-5	9



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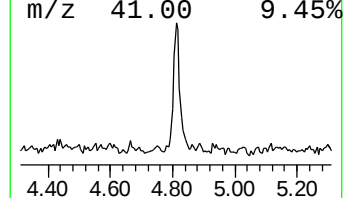
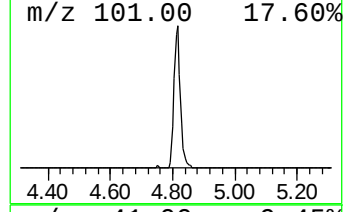
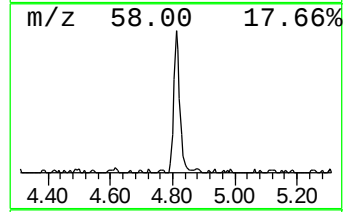
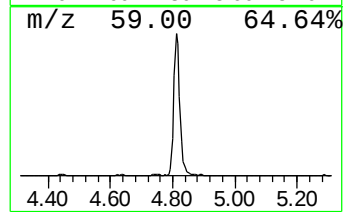
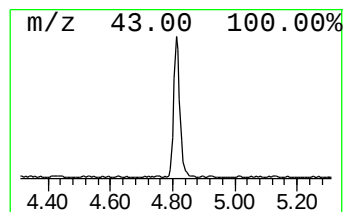
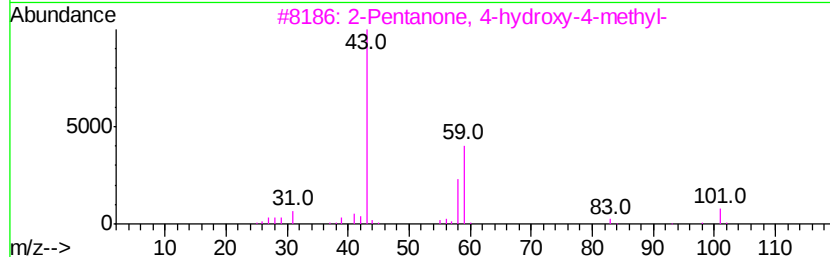
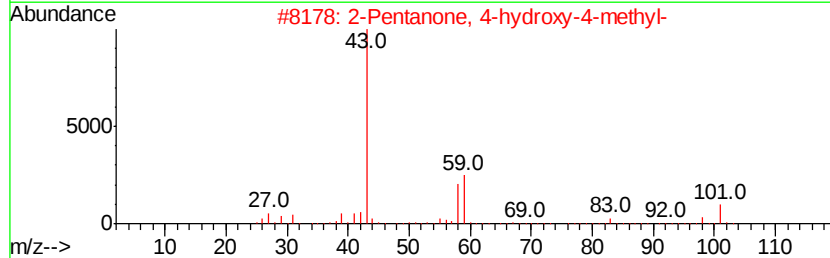
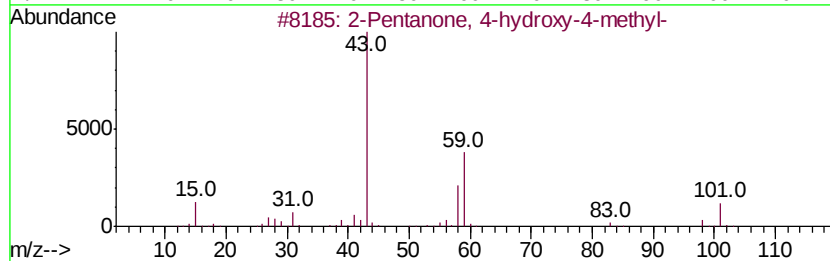
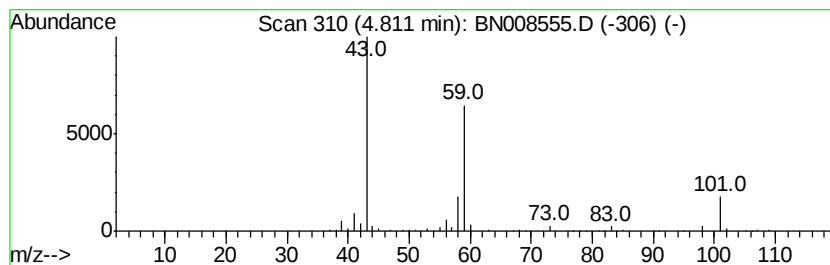
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	3.88 ng/ul	399856	1,4-Dichlorobenzene-d4	7.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
4			Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	40
5			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33



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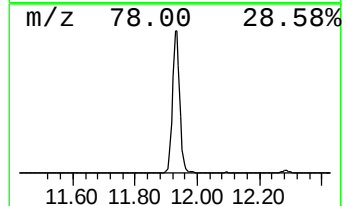
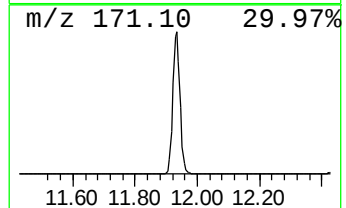
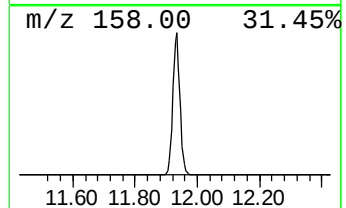
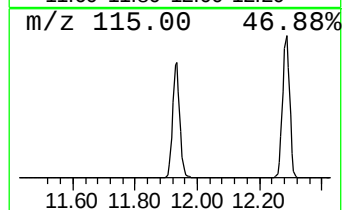
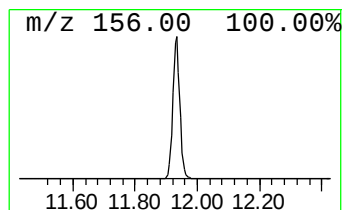
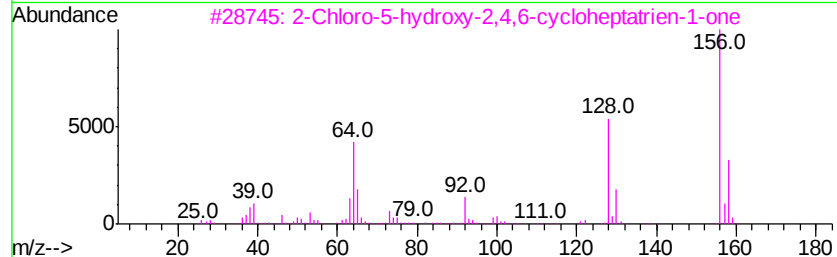
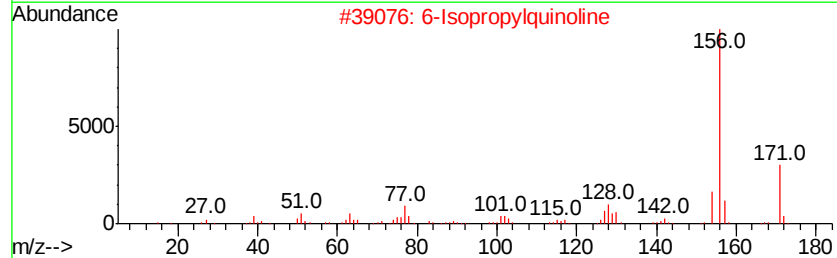
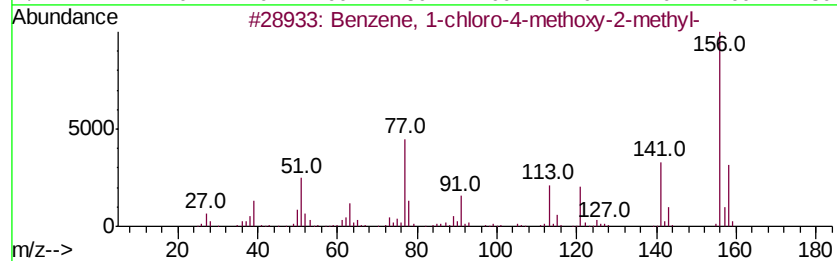
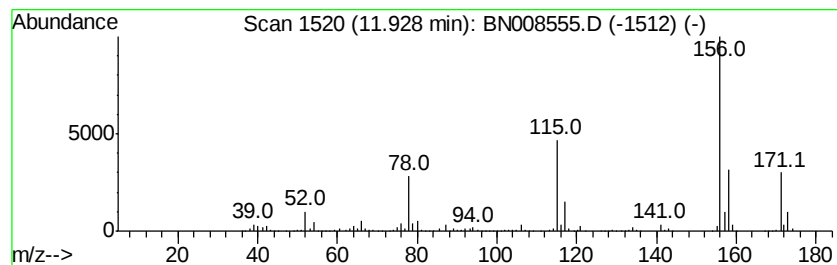
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	36.85 ng/ul	5644970	Napthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	38
2		6-Isopropylquinoline	171	C12H13N	000135-79-5	38
3		2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	32
4		Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32
5		Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28



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Data File : BN008555.D
Acq On : 14 Nov 2019 13:53
Operator : JU
Sample : K5656-10
Misc :
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Instrument :
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ClientSampleId :
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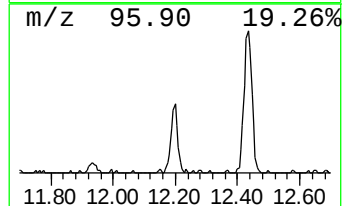
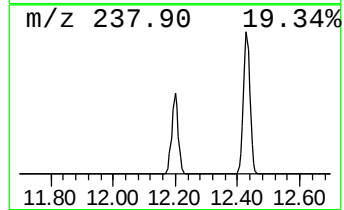
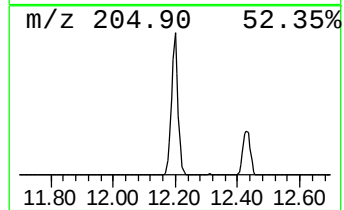
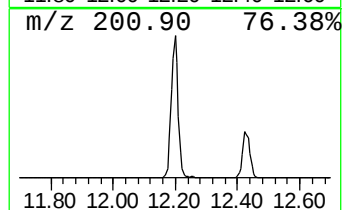
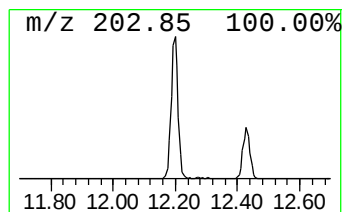
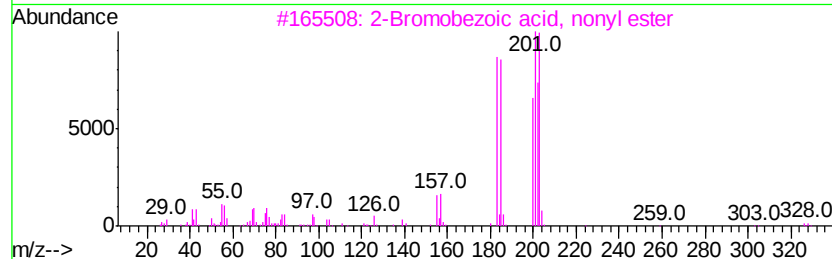
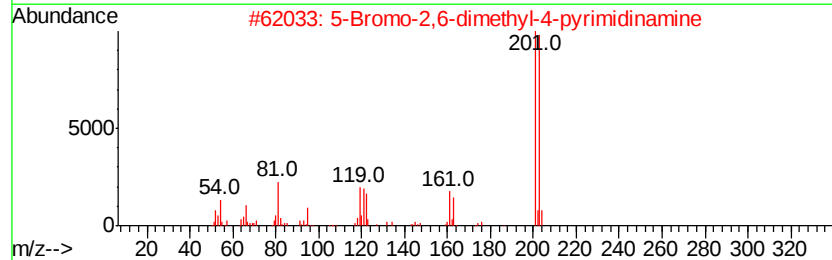
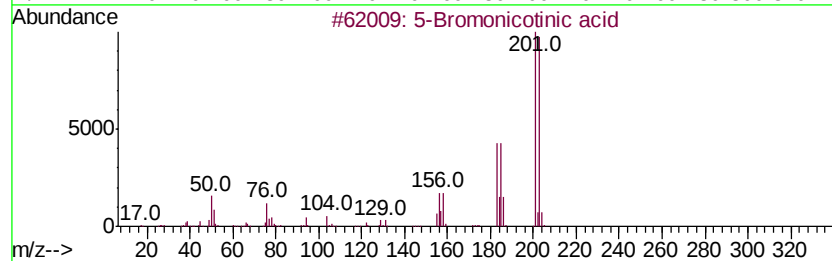
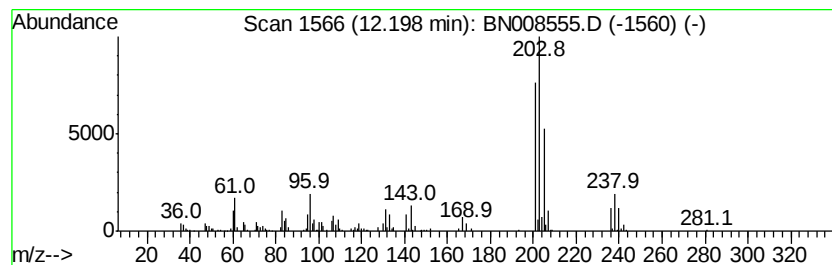
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.20	3.52 ng/ul	539473	Naphthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Bromonicotinic acid	201	C6H4BrNO2	020826-04-4	43
2		5-Bromo-2,6-dimethyl-4-pyrimidin...	201	C6H8BrN3	007752-80-9	43
3		2-Bromobezoic acid, nonyl ester	326	C16H23BrO2	1000282-80-5	43
4		1-Benzenecarbonyl chloride, 2,4-...	238	C8H5Cl3O2	1000196-85-1	43
5		4-Amino-5,7-dichlorobenzofurazan	203	C6H3Cl2N3O	330982-41-7	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
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ALS Vial : 32 Sample Multiplier: 1

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ClientSampleId :
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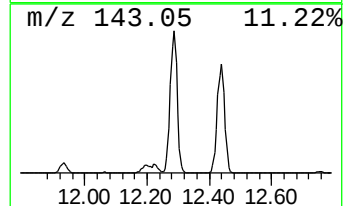
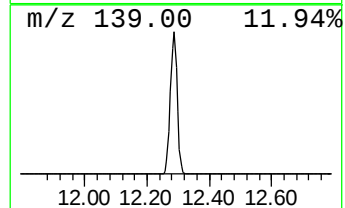
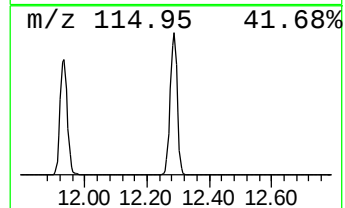
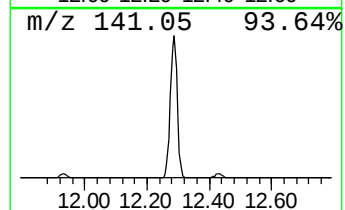
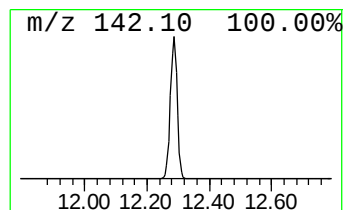
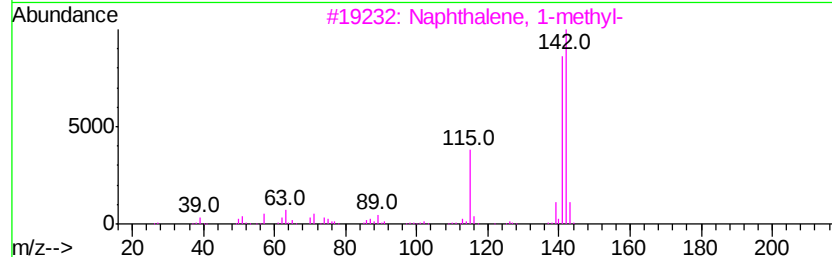
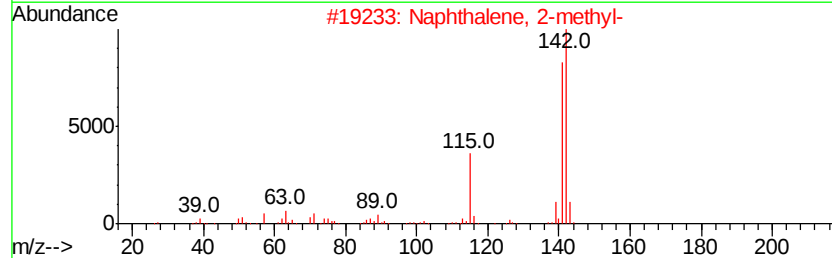
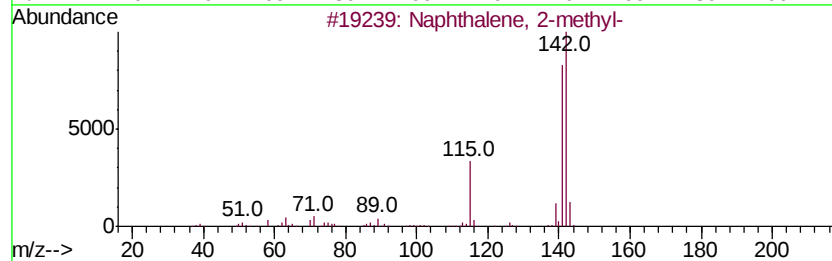
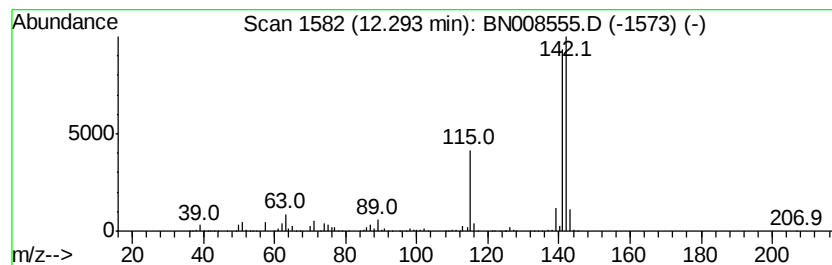
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	44.86 ng/ul	6871410	Naphthalene-d8	10.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008555.D
Acq On : 14 Nov 2019 13:53
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Sample : K5656-10
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ClientSampleId :
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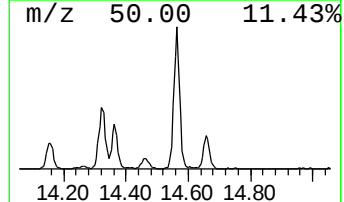
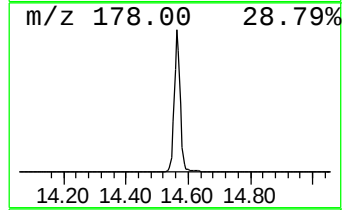
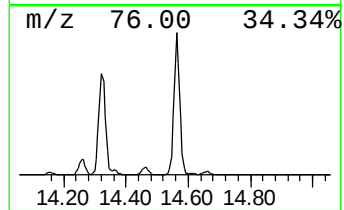
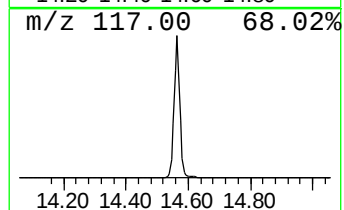
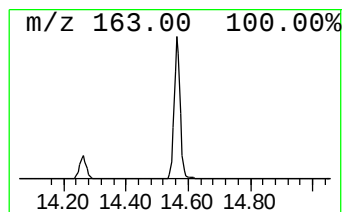
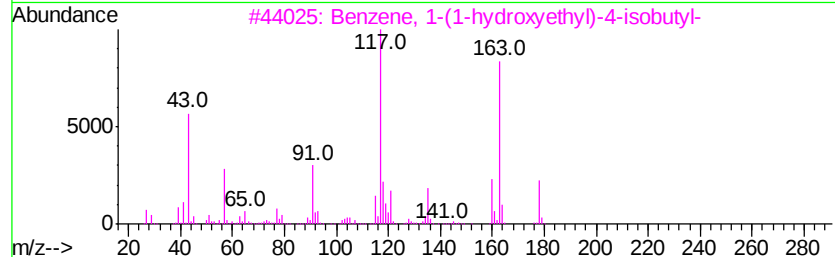
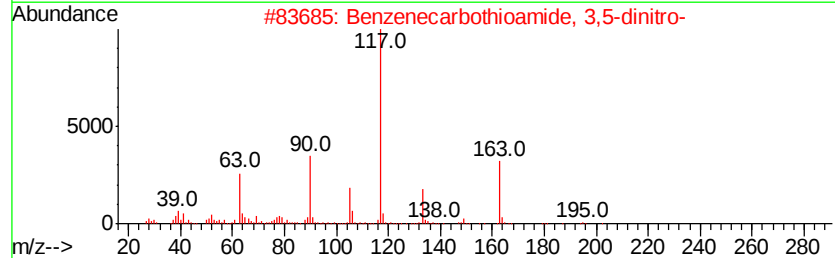
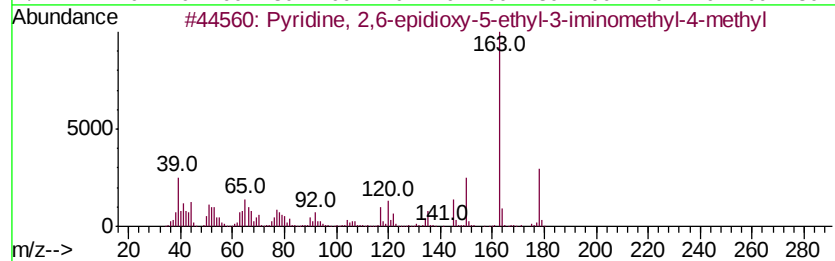
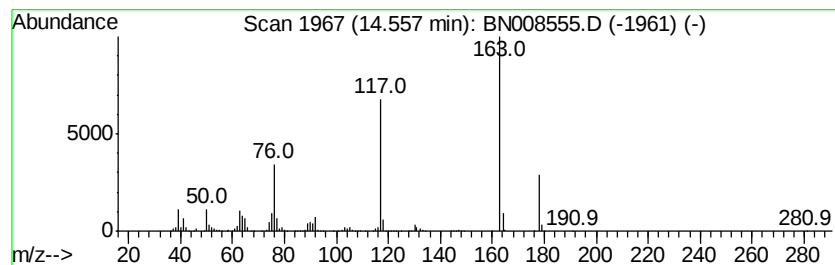
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Pyridine, 2,6-epidioxv-5-et... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	21.37 ng/ul	4131970	Acenaphthene-d10	14.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyridine, 2,6-epidioxv-5-ethyl-3...	178	C9H10N2O2	1000263-31-1	53
2		Benzenecarbothioamide, 3,5-dinitro-	227	C7H5N3O4S	037773-67-4	53
3		Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	50
4		Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	50
5		Methyl p-chlorophenylethynyl ketone	178	C10H7ClO	039180-26-2	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008555.D
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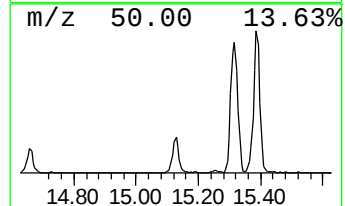
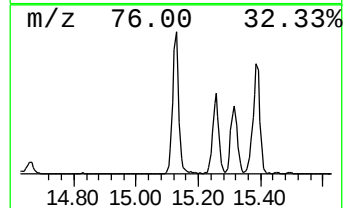
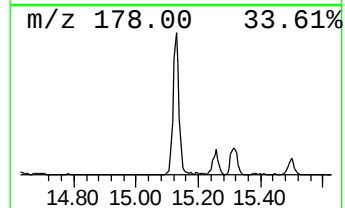
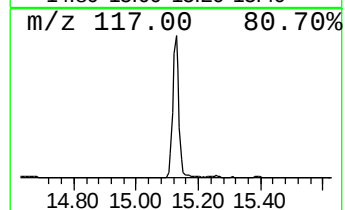
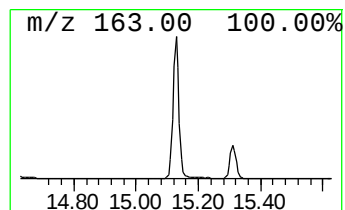
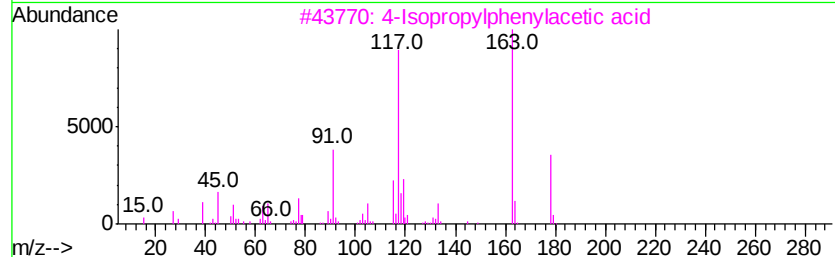
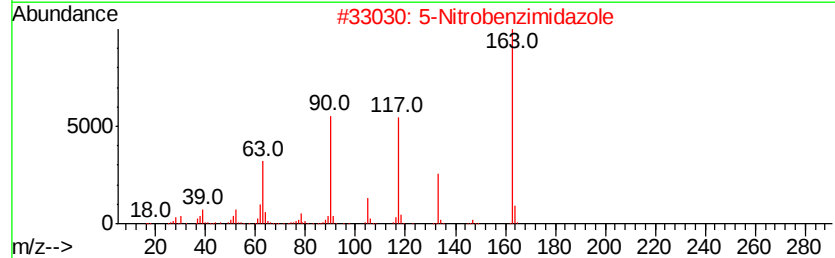
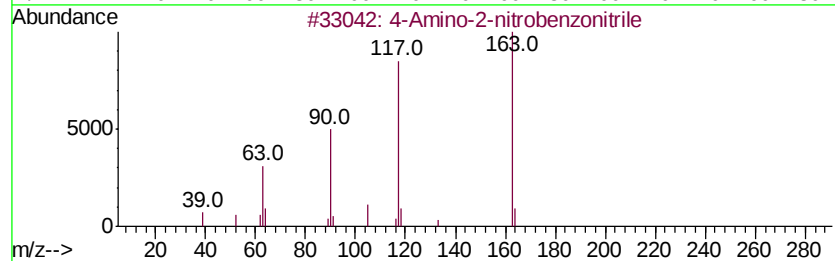
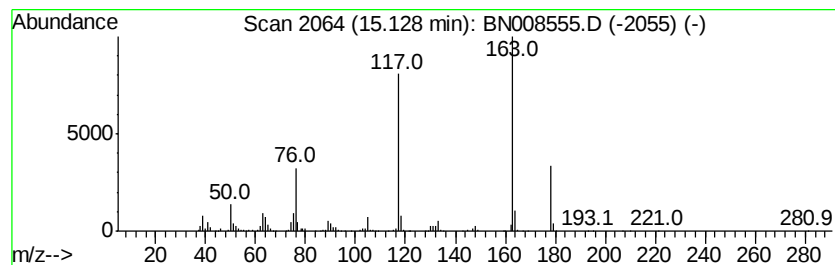
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	6.49 ng/ul	1255090	Acenaphthene-d10	14.26
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	4-Amino-2-nitrobenzonitrile	163 C7H5N3O2	072115-07-2	64
2	5-Nitrobenzimidazole	163 C7H5N3O2	1000229-89-5	59
3	4-Isopropylphenylacetic acid	178 C11H14O2	004476-28-2	45
4	Propofol	178 C12H18O	002078-54-8	43
5	Propofol	178 C12H18O	002078-54-8	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008555.D
Acq On : 14 Nov 2019 13:53
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Misc :
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Instrument :
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ClientSampleId :
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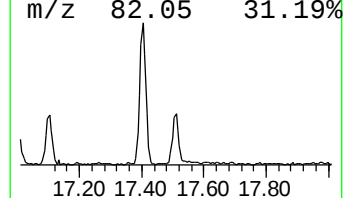
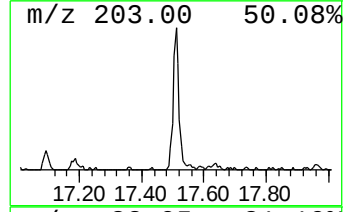
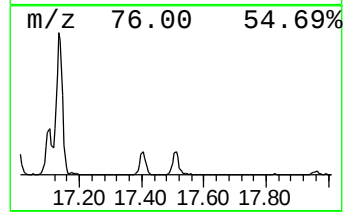
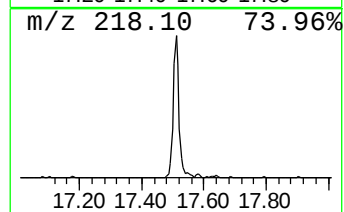
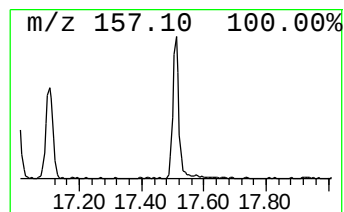
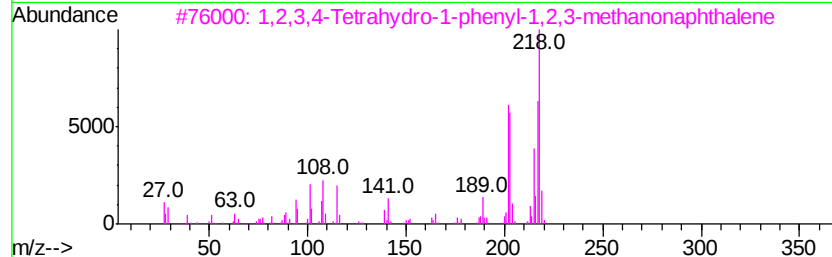
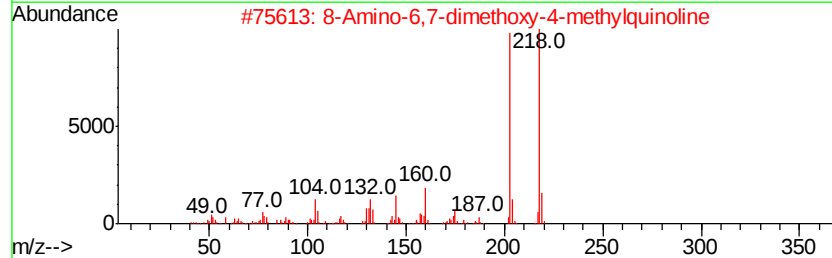
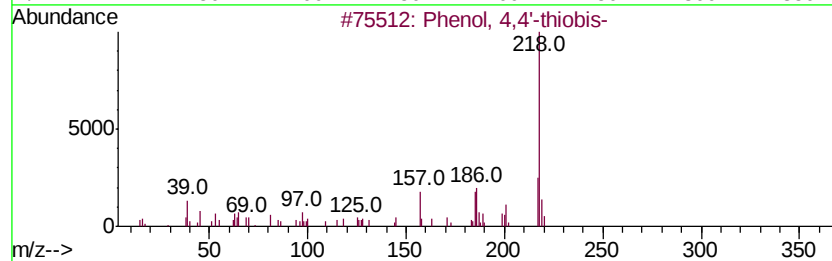
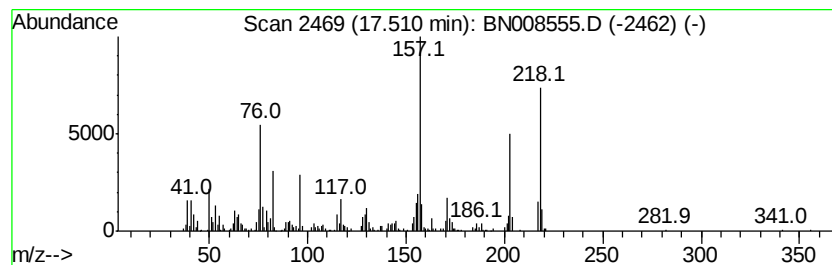
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.51	2.58 ng/ul	576199	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-thiobis-	218	C12H10O2S	002664-63-3	38
2		8-Amino-6,7-dimethoxy-4-methylqu...	218	C12H14N2O2	1000213-07-2	30
3		1,2,3,4-Tetrahydro-1-phenyl-1,2,...	218	C17H14	138089-69-7	30
4		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	30
5		Benzenamine, 2,6-dimethyl-N-(4,4...	218	C13H18N2O	281211-68-5	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008555.D
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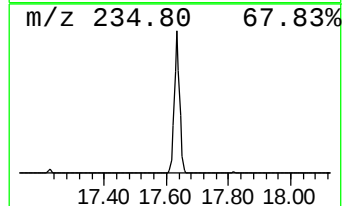
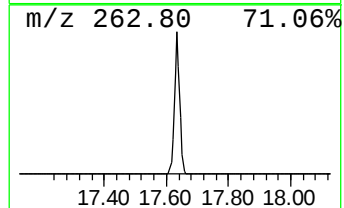
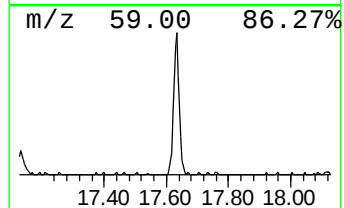
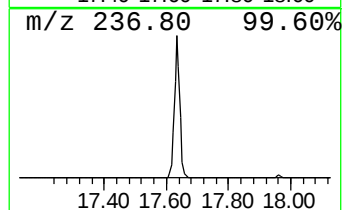
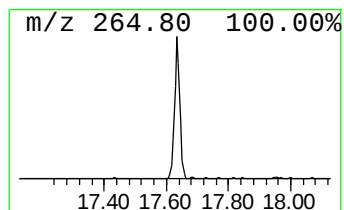
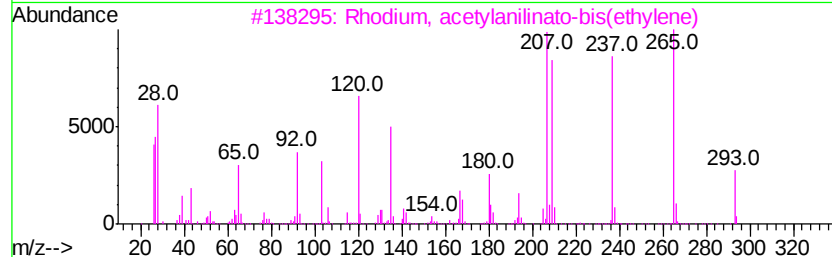
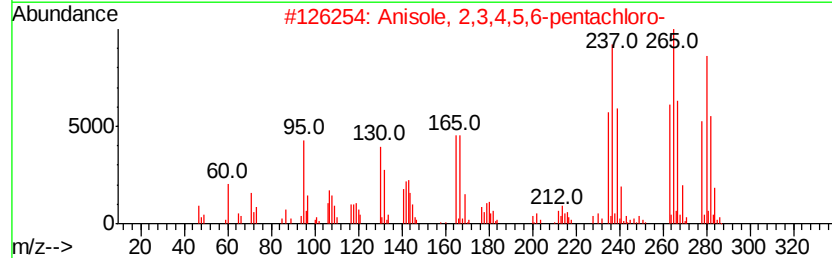
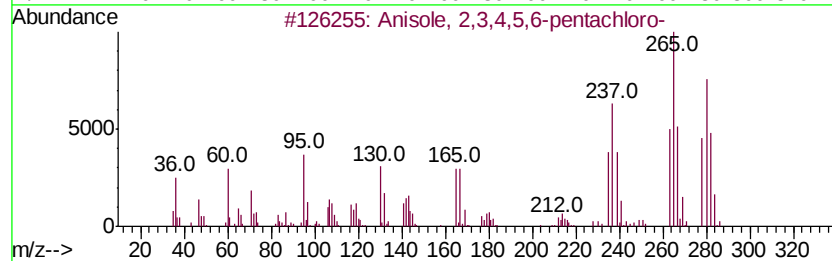
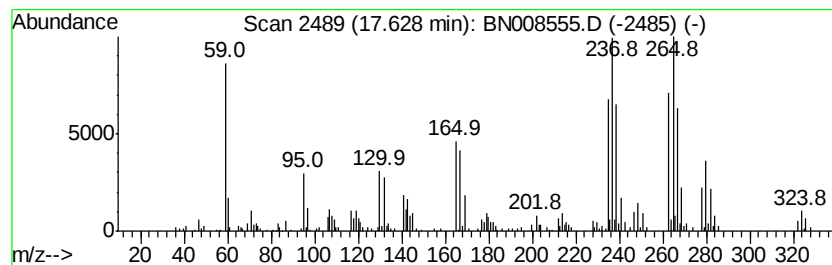
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Anisole, 2,3,4,5,6-pentachl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.63	2.65 ng/ul	591702	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	60
2		Anisole, 2,3,4,5,6-pentachloro-	278	C7H3Cl5O	001825-21-4	46
3		Rhodium, acetylanilinato-bis(eth...	293	C12H16N0Rh	1000153-76-6	30
4		2,3,4,5,6-Pentachloroaniline	263	C6H2Cl5N	000527-20-8	30
5		Hexachlorocyclopentadiene	270	C5Cl6	000077-47-4	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008555.D
Acq On : 14 Nov 2019 13:53
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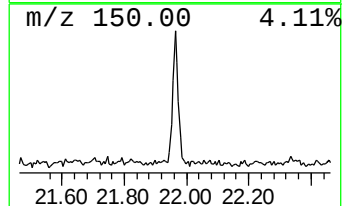
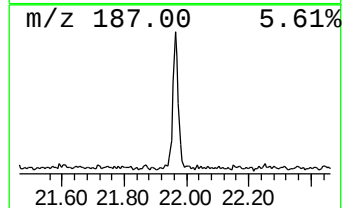
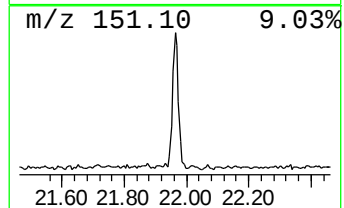
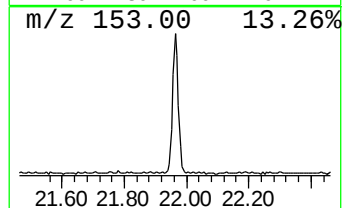
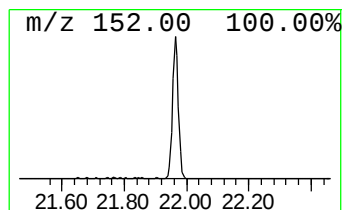
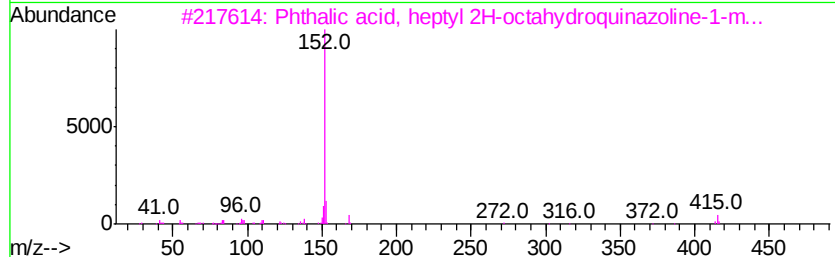
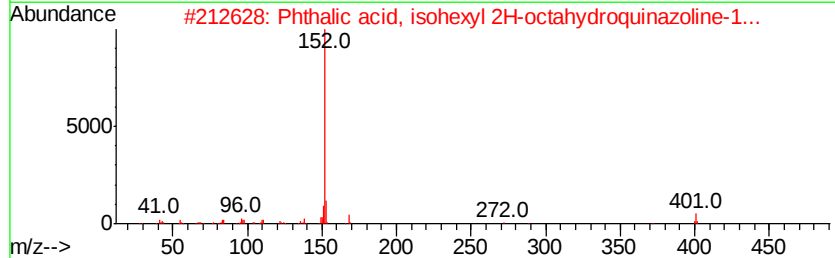
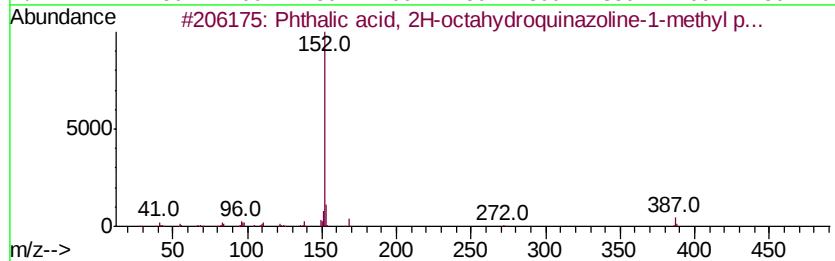
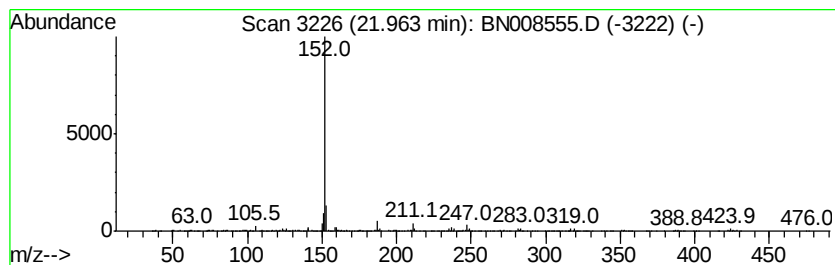
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phthalic acid, 2H-octahydro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.96	5.40 ng/ul	1271800	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, 2H-octahydroquina...	387	C23H33N04	1000322-80-5	72
2		Phthalic acid, isohexyl 2H-octah...	401	C24H35N04	1000322-80-6	72
3		Phthalic acid, heptyl 2H-octahvd...	415	C25H37N04	1000322-80-8	72
4		Cyclobuta(1,2-a:3,4-a')diacenaph...	304	C24H16	007259-03-2	72
5		Acenaphthylene	152	C12H8	000208-96-8	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN111319\
Data File : BN008555.D
Acq On : 14 Nov 2019 13:53
Operator : JU
Sample : K5656-10
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A5458

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.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
Glycolaldehyde di...	3.02	2.6	ng/ul	272324	1	7.63	2062600	20.0
2-Pentanone, 4-hy...	4.81	3.9	ng/ul	399856	1	7.63	2062600	20.0
unknown-01	11.93	36.9	ng/ul	5644970	2	10.40	3063510	20.0
unknown-02	12.20	3.5	ng/ul	539473	2	10.40	3063510	20.0
Naphthalene, 1-me...	12.29	44.9	ng/ul	6871410	2	10.40	3063510	20.0
Pyridine, 2,6-epi...	14.56	21.4	ng/ul	4131970	3	14.26	3867050	20.0
4-Amino-2-nitrobe...	15.13	6.5	ng/ul	1255090	3	14.26	3867050	20.0
unknown-03	17.51	2.6	ng/ul	576199	4	17.00	4461170	20.0
Anisole, 2,3,4,5,...	17.63	2.6	ng/ul	591702	4	17.00	4461170	20.0
Phthalic acid, 2H...	21.96	5.4	ng/ul	1271800	5	21.20	4709370	20.0