

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF011719\
 Data File : BF112089.D
 Acq On : 17 Jan 2019 15:11
 Operator : JU/SJ
 Sample : K1167-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D-9

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF011619.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.152	6	11	19	rVB	192503	265742	2.88%	0.307%
2	3.398	221	223	243	rBV	60318	178749	1.94%	0.206%
3	5.087	506	510	518	rVB	372724	527651	5.71%	0.609%
4	5.504	576	581	584	rBV	7299637	7700257	83.36%	8.885%
5	5.704	609	615	625	rBV	144106	174626	1.89%	0.201%
6	6.510	747	752	761	rBV2	6877250	8849163	95.80%	10.211%
7	6.639	769	774	777	rBV	8463170	8404019	90.98%	9.697%
8	6.698	780	784	788	rVV2	222611	272531	2.95%	0.314%
9	6.739	788	791	796	rVB	111490	107967	1.17%	0.125%
10	6.863	808	812	819	rBV	2203408	1893484	20.50%	2.185%
11	7.022	834	839	842	rBV	7361798	6605464	71.51%	7.622%
12	7.122	852	856	863	rBV	143565	151319	1.64%	0.175%
13	7.422	901	907	911	rBV	4765656	5379881	58.24%	6.208%
14	7.716	954	957	968	rVB	405624	367402	3.98%	0.424%
15	8.145	1026	1030	1033	rBV	2580711	2387908	25.85%	2.755%
16	8.357	1061	1066	1069	rBV	347712	285828	3.09%	0.330%
17	8.416	1073	1076	1080	rBV	403408	331525	3.59%	0.383%
18	8.851	1145	1150	1155	rVB2	149622	182830	1.98%	0.211%
19	9.133	1196	1198	1201	rBV	106434	95854	1.04%	0.111%
20	9.222	1208	1213	1216	rBV	9310236	9236865	100.00%	10.658%
21	9.345	1231	1234	1237	rBV	119552	131017	1.42%	0.151%
22	9.604	1275	1278	1283	rBV	1057553	975101	10.56%	1.125%
23	9.757	1301	1304	1309	rBV	120480	123408	1.34%	0.142%
24	9.898	1324	1328	1331	rBV2	3018272	2512265	27.20%	2.899%
25	9.927	1331	1333	1337	rVB	191674	164661	1.78%	0.190%
26	10.445	1418	1421	1426	rVB	185762	179620	1.94%	0.207%
27	10.692	1458	1463	1466	rBV	5805214	5545119	60.03%	6.398%
28	10.816	1479	1484	1490	rVB	151755	192887	2.09%	0.223%
29	11.280	1560	1563	1567	rVB	123623	118941	1.29%	0.137%
30	11.386	1577	1581	1583	rBV	2583655	2364723	25.60%	2.729%
31	11.410	1583	1585	1589	rVV	713811	599098	6.49%	0.691%
32	11.457	1591	1593	1596	rVB	208903	172811	1.87%	0.199%
33	11.886	1663	1666	1668	rBV	106261	101351	1.10%	0.117%
34	11.910	1668	1670	1676	rVB2	317175	297306	3.22%	0.343%

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Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF011619.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.998	1682	1685	1688	rBV	225568	239311	2.59%	0.276%
36	12.598	1784	1787	1791	rBV	1261502	1017150	11.01%	1.174%
37	12.827	1820	1826	1832	rBV	1077261	1123108	12.16%	1.296%
38	12.974	1846	1851	1854	rBV	7780961	7798629	84.43%	8.999%
39	13.157	1880	1882	1886	rVB	124023	105727	1.14%	0.122%
40	13.663	1965	1968	1972	rVB	86896	101166	1.10%	0.117%
41	13.857	1998	2001	2008	rVB2	787226	876512	9.49%	1.011%
42	14.027	2025	2030	2032	rBV2	2463372	2547282	27.58%	2.939%
43	14.051	2032	2034	2039	rVB	504808	426081	4.61%	0.492%
44	14.180	2054	2056	2060	rVB	118352	108965	1.18%	0.126%
45	14.486	2105	2108	2115	rBV3	270437	372469	4.03%	0.430%
46	14.792	2157	2160	2165	rVB	236477	254674	2.76%	0.294%
47	15.068	2203	2207	2210	rBV	421230	591966	6.41%	0.683%
48	15.133	2215	2218	2222	rVV	293786	319043	3.45%	0.368%
49	15.374	2256	2259	2265	rVB	260116	344318	3.73%	0.397%
50	15.439	2266	2270	2275	rVV	292473	403152	4.36%	0.465%
51	15.504	2276	2281	2293	rVB	1542639	2281349	24.70%	2.632%
52	15.951	2354	2357	2363	rVB	214223	286674	3.10%	0.331%
53	16.462	2440	2444	2447	rBV	113215	164415	1.78%	0.190%
54	16.980	2528	2532	2539	rVB2	126087	225848	2.45%	0.261%
55	17.421	2604	2607	2615	rVB	102675	200132	2.17%	0.231%

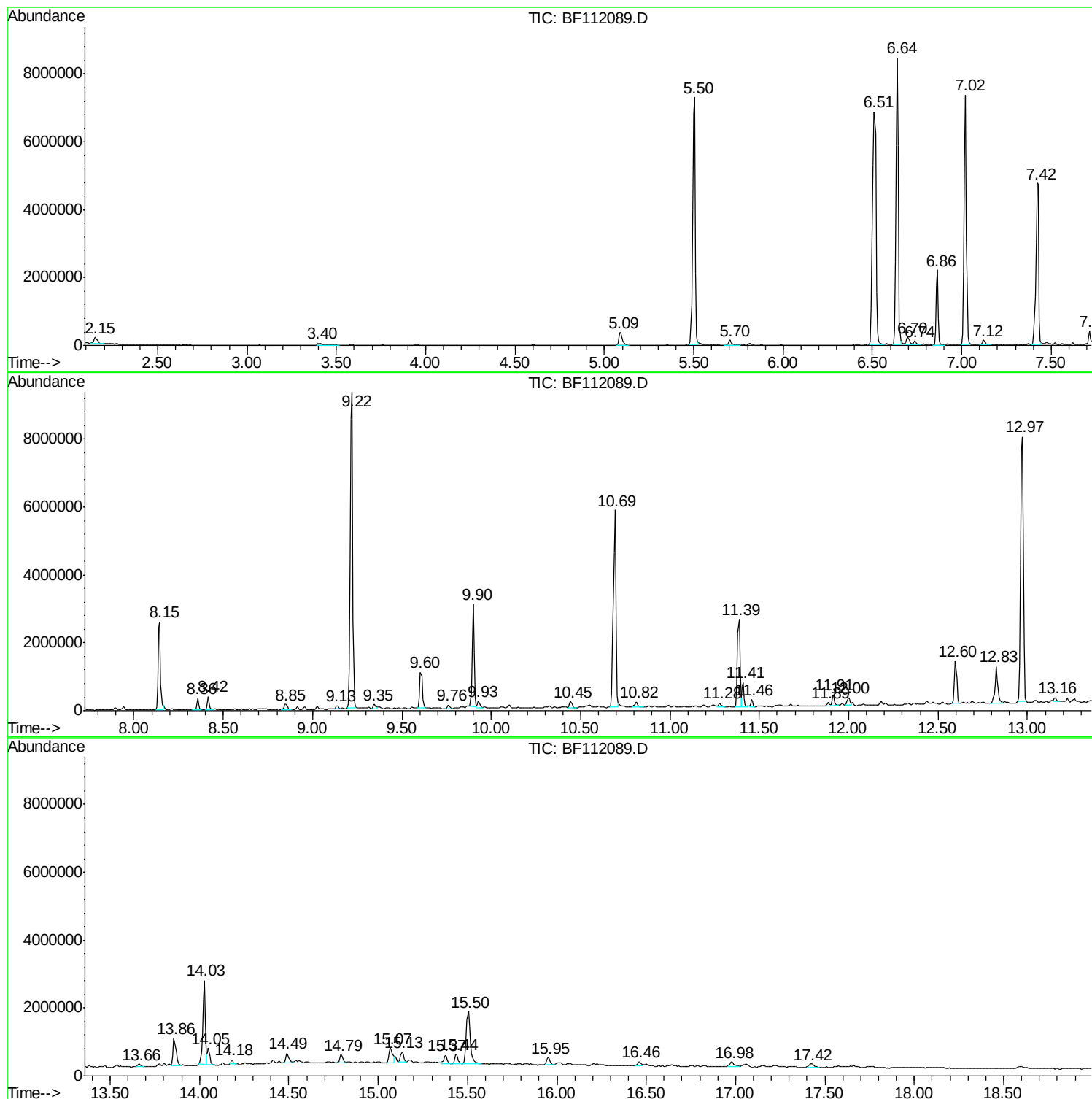
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF011619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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ClientSampleId :
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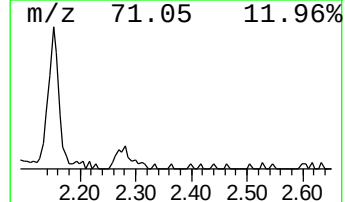
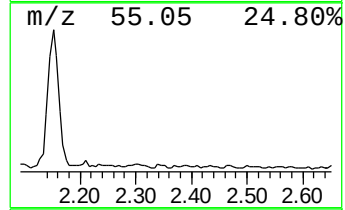
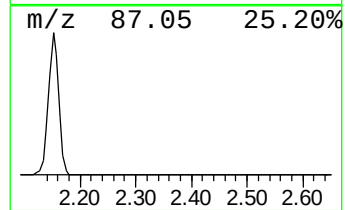
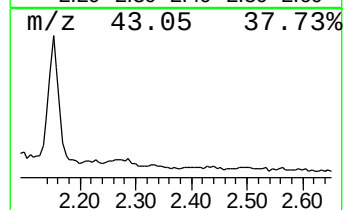
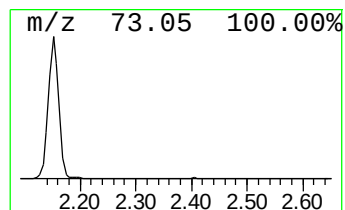
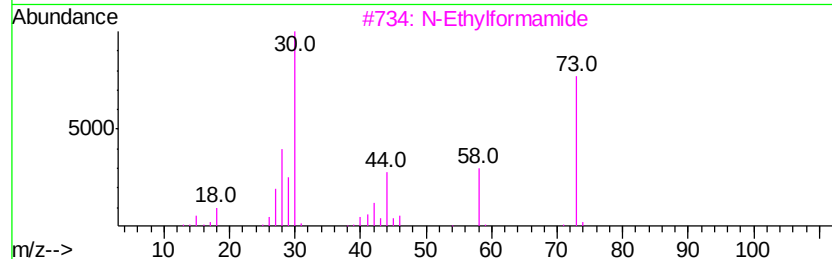
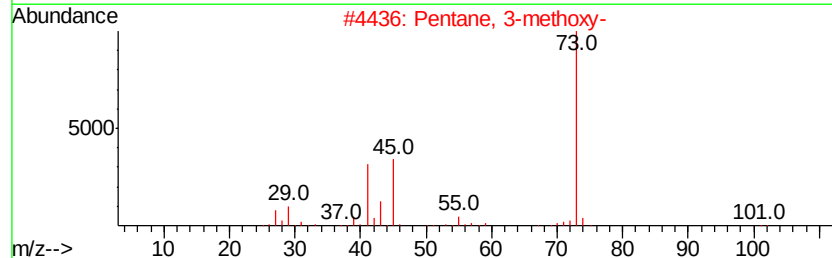
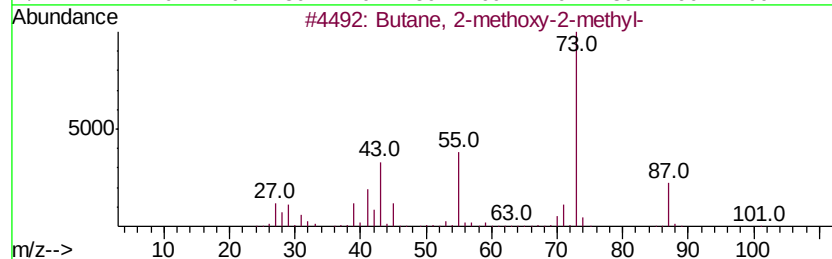
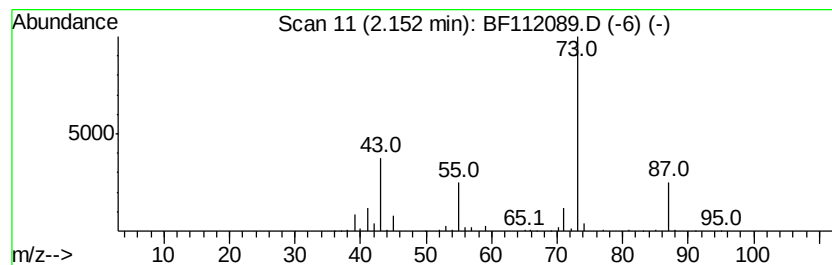
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	2.81 ng	265742	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
3			N-Ethylformamide	73	C3H7NO	000627-45-2	9
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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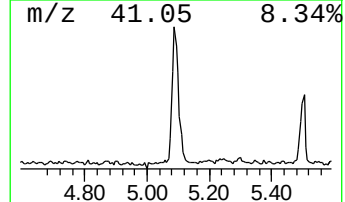
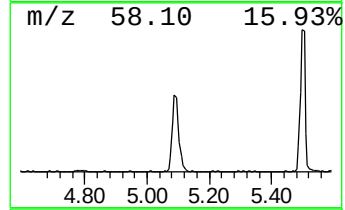
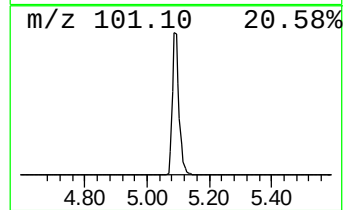
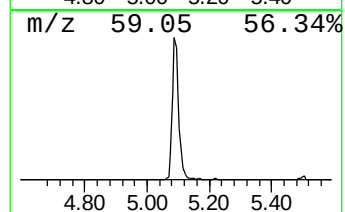
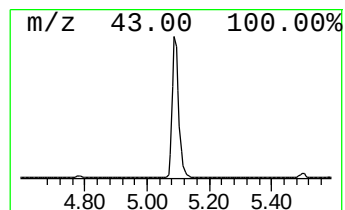
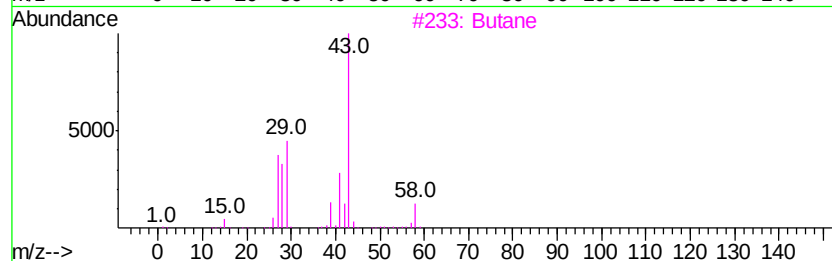
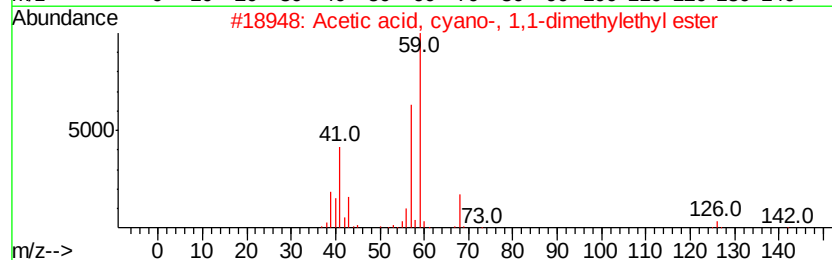
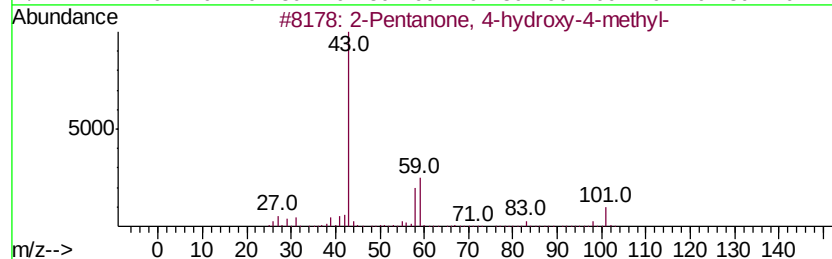
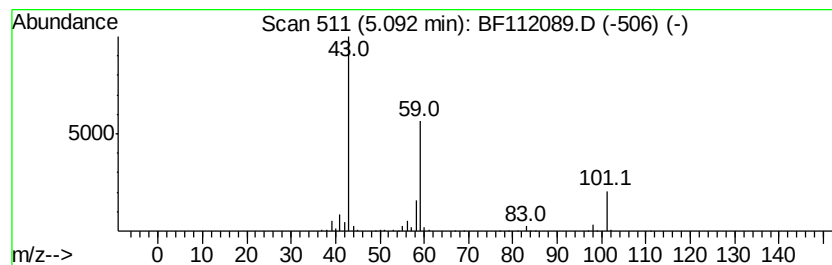
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.09	5.57 ng	527651	1,4-Dichlorobenzene-d4	6.86	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3	Butane	58	C4H10	000106-97-8	9
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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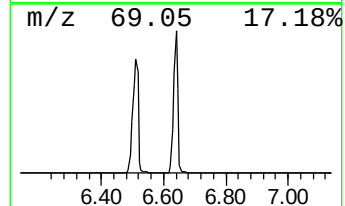
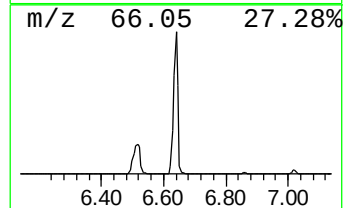
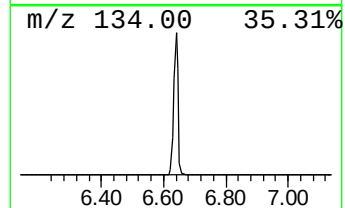
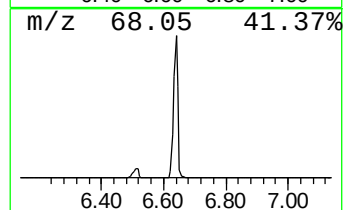
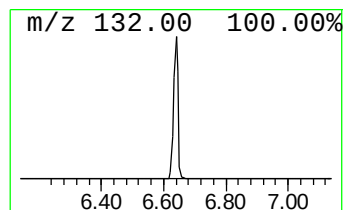
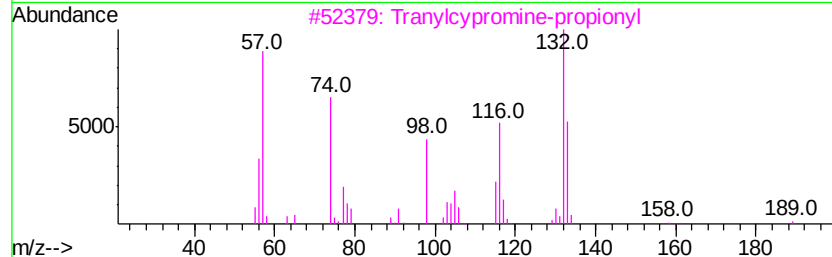
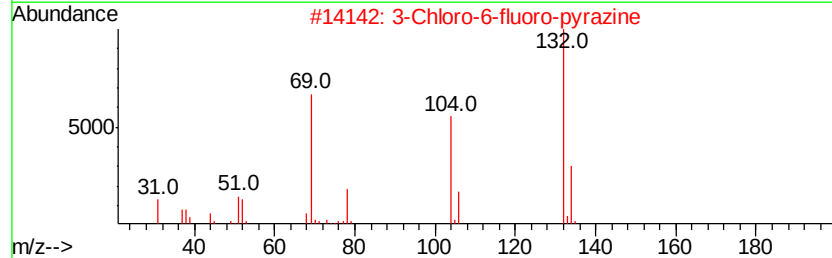
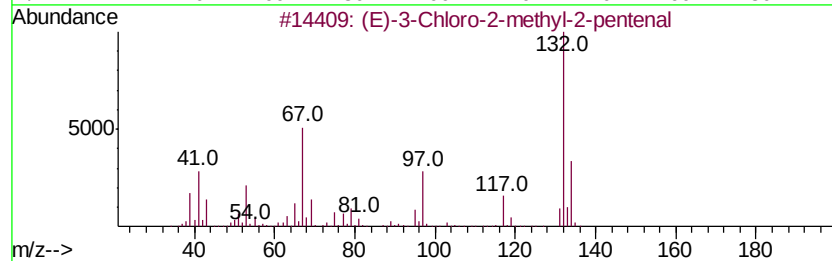
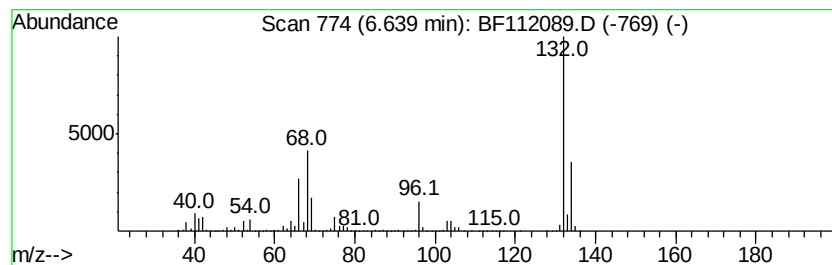
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	88.77 ng	8404020	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	30
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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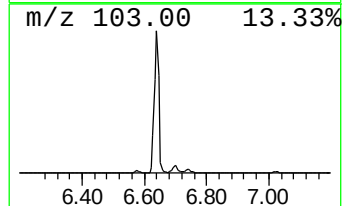
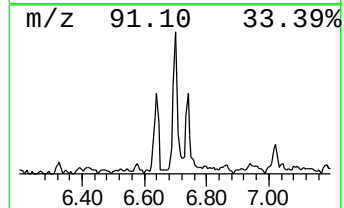
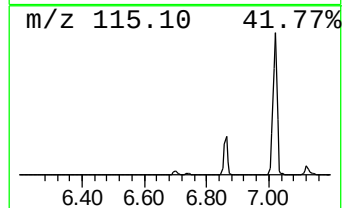
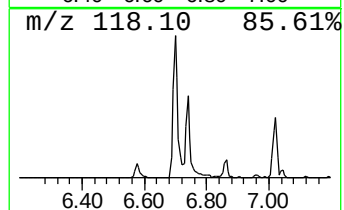
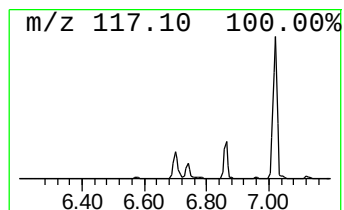
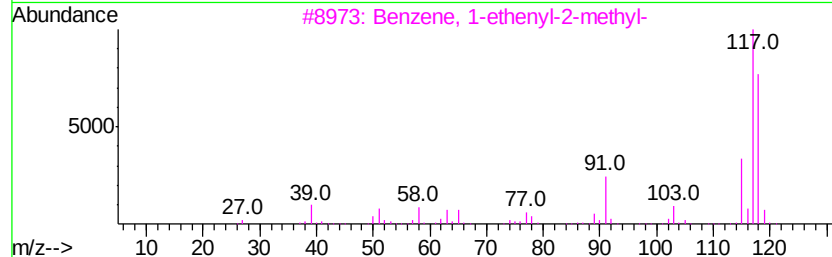
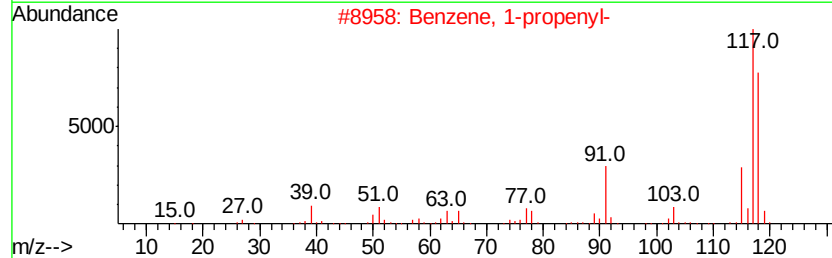
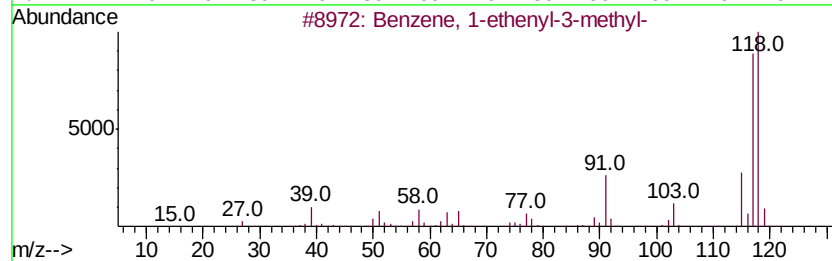
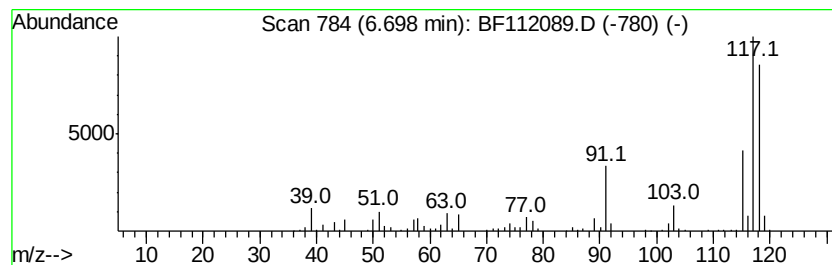
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, 1-ethenyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	2.88 ng	272531	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	94
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
4			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	93
5			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	91



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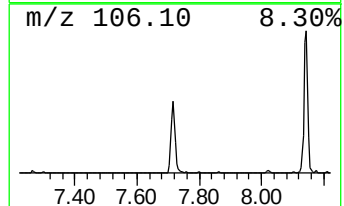
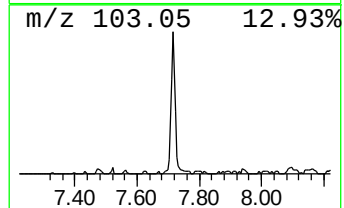
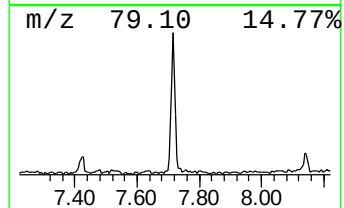
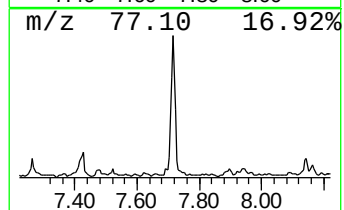
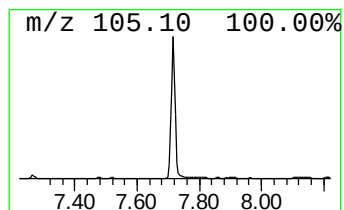
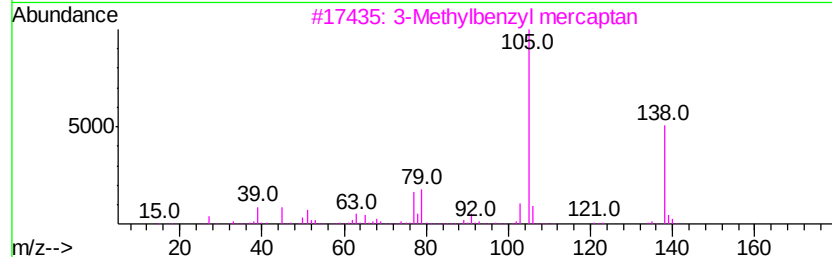
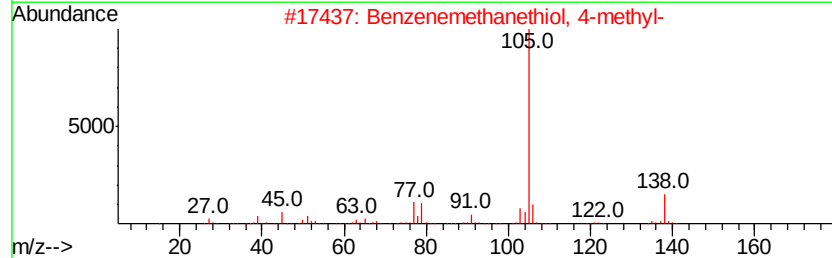
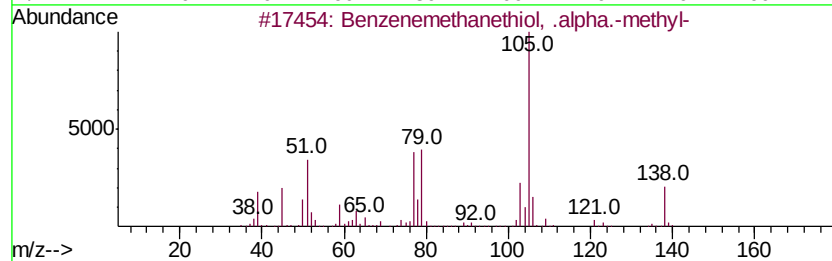
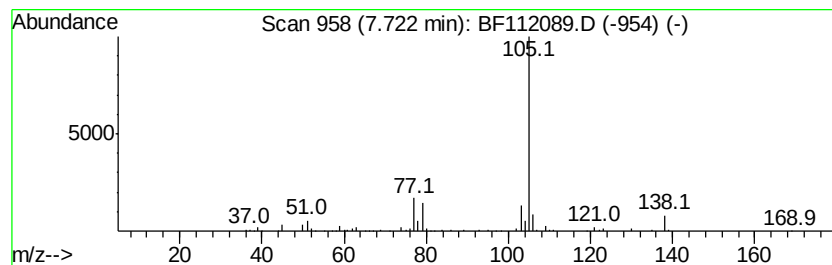
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzenemethanethiol, .alpha... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	3.08 ng	367402	Napthalene-d8	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenemethanethiol, .alpha.-met...	138	C8H10S	006263-65-6	86
2		Benzenemethanethiol, 4-methyl-	138	C8H10S	004498-99-1	80
3		3-Methylbenzyl mercaptan	138	C8H10S	025697-56-7	72
4		Benzene, (1-bromoethyl)-	184	C8H9Br	000585-71-7	72
5		3,4-Dimethylthiophenol	138	C8H10S	018800-53-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
Data File : BF112089.D
Acq On : 17 Jan 2019 15:11
Operator : JU/SJ
Sample : K1167-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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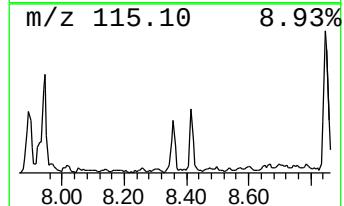
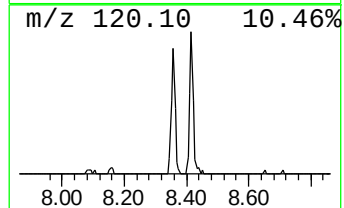
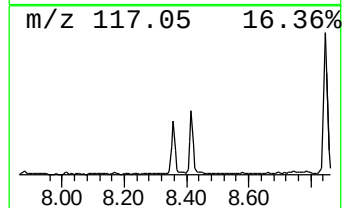
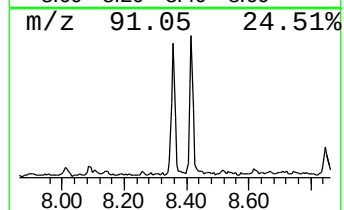
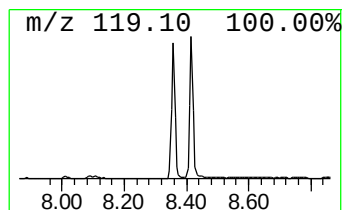
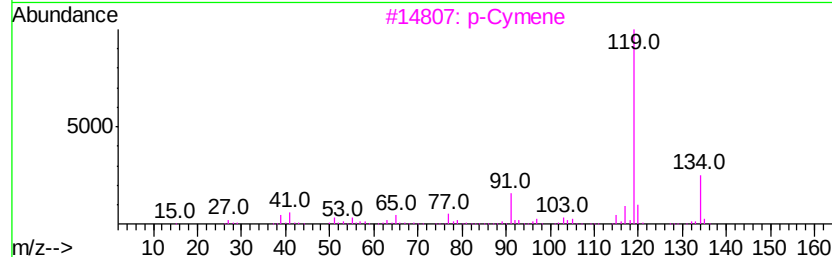
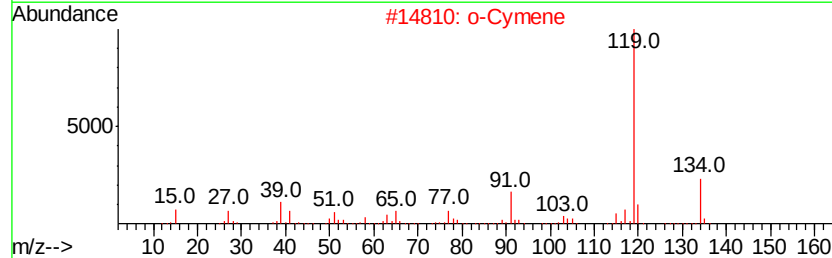
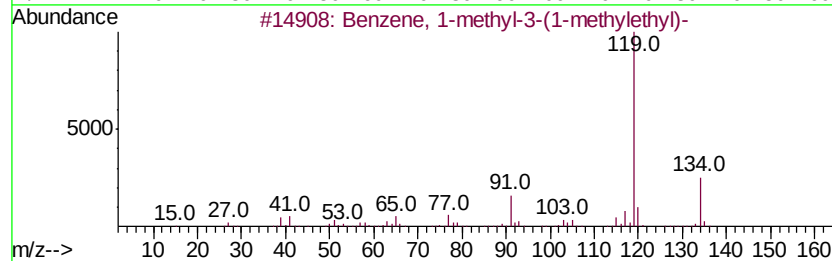
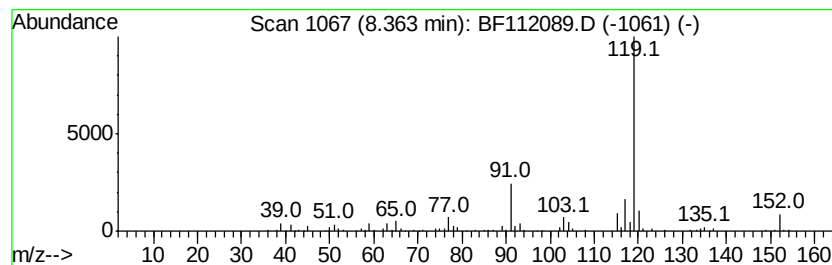
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1-methyl-3-(1-meth... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.36	2.39 ng	285828	Napthalene-d8	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	83
2		o-Cymene	134	C10H14	000527-84-4	81
3		p-Cymene	134	C10H14	000099-87-6	80
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	72
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
Data File : BF112089.D
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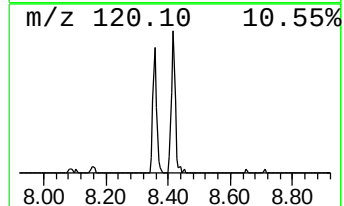
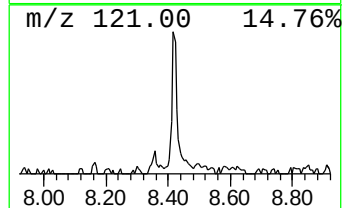
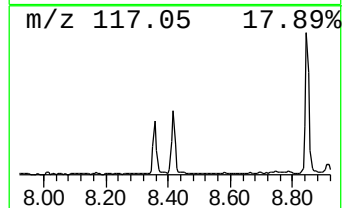
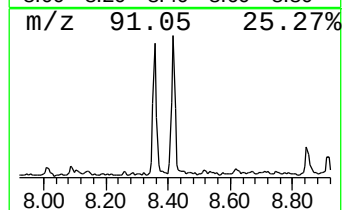
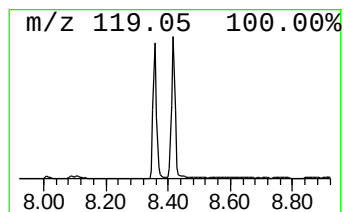
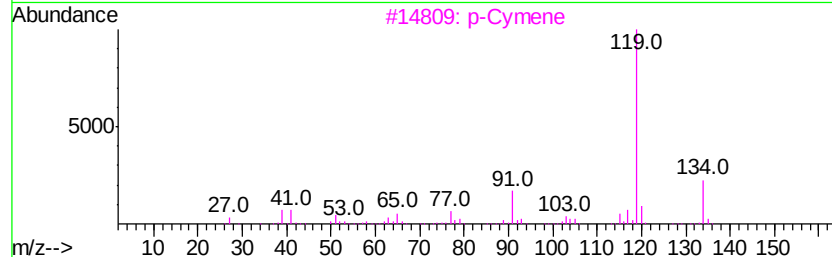
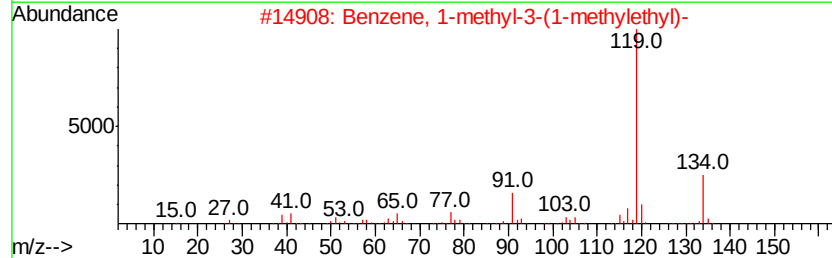
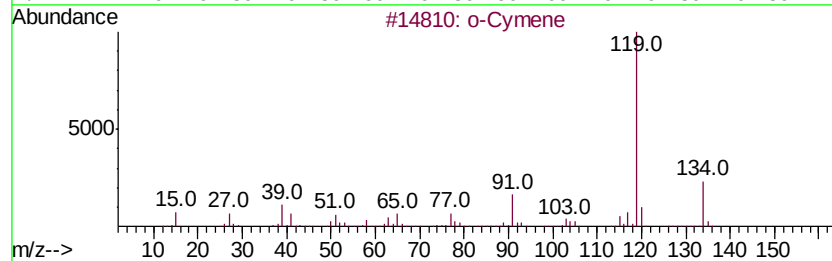
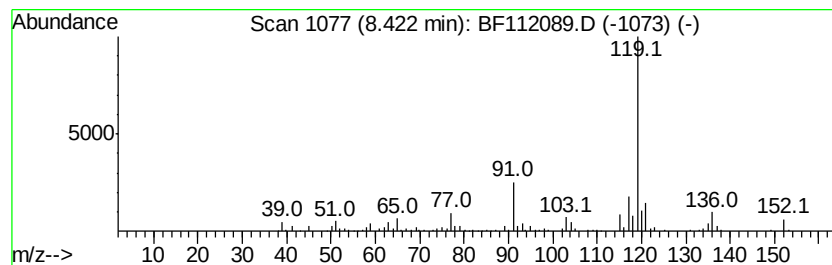
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	2.78 ng	331525	Napthalene-d8	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	81
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	81
3		p-Cymene	134	C10H14	000099-87-6	74
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	72
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
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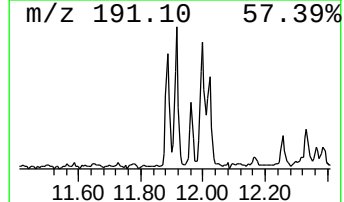
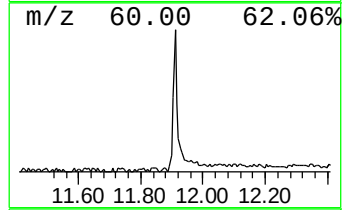
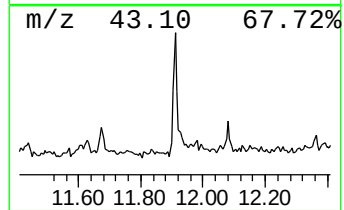
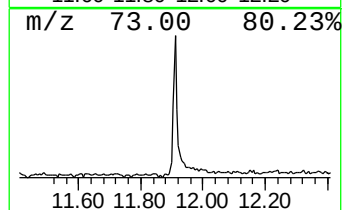
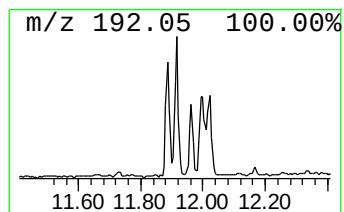
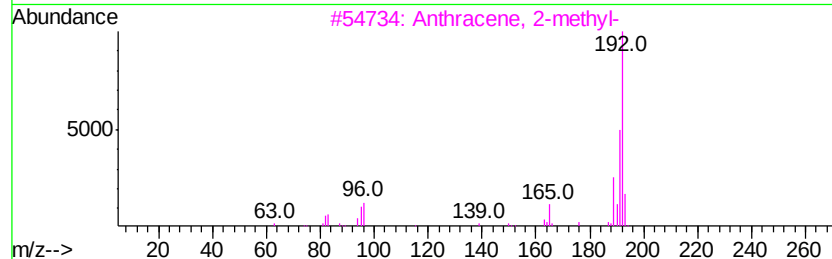
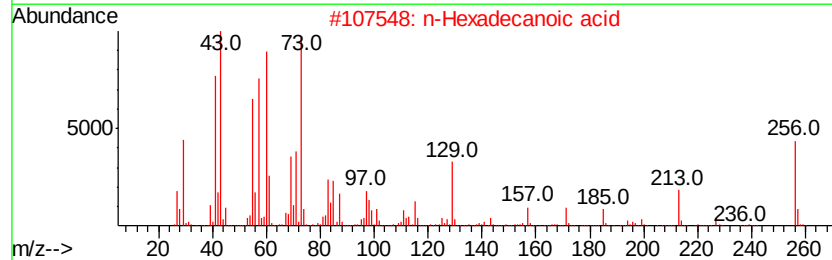
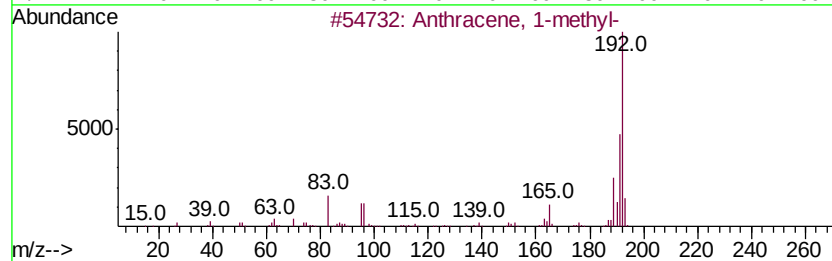
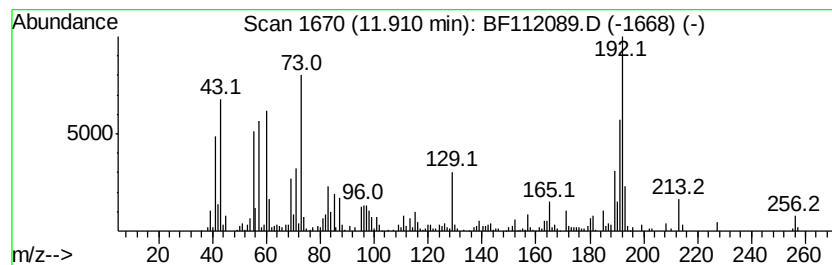
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Anthracene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.91	2.51 ng	297306	Phenanthrene-d10		11.39
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
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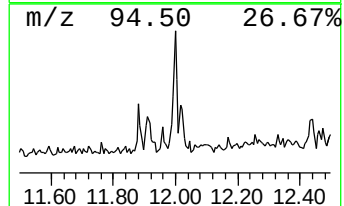
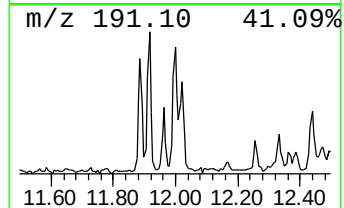
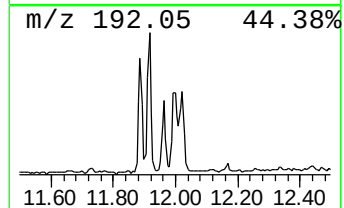
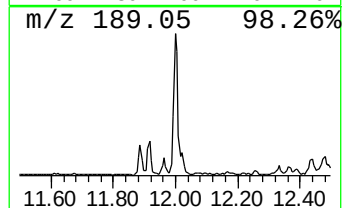
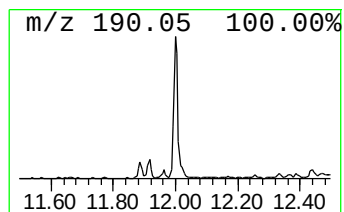
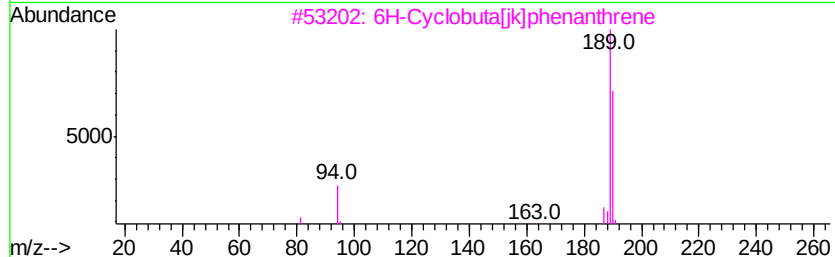
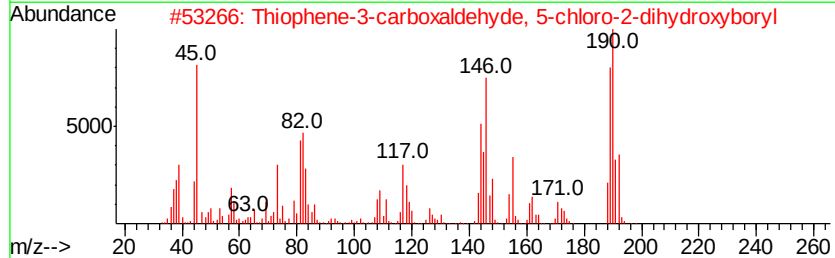
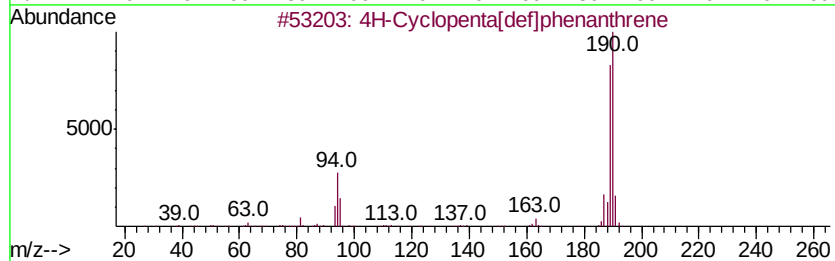
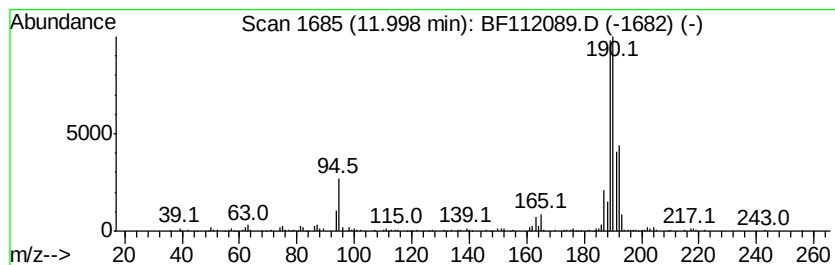
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	2.02 ng	239311	Phenanthrene-d10	11.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
3			6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	53
4			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	43
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
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Instrument :
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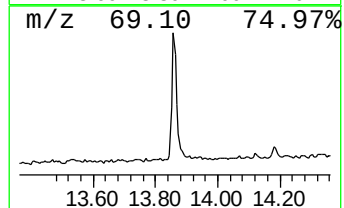
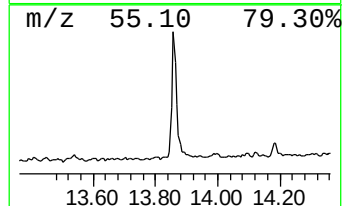
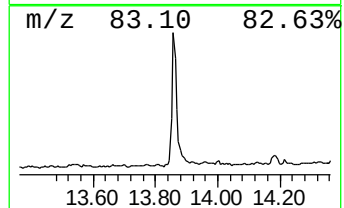
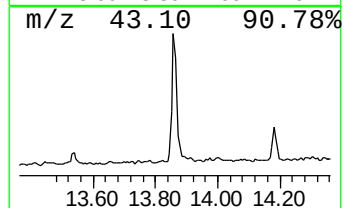
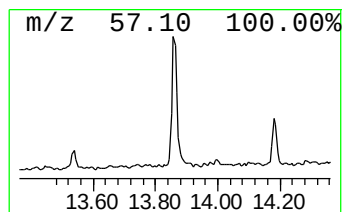
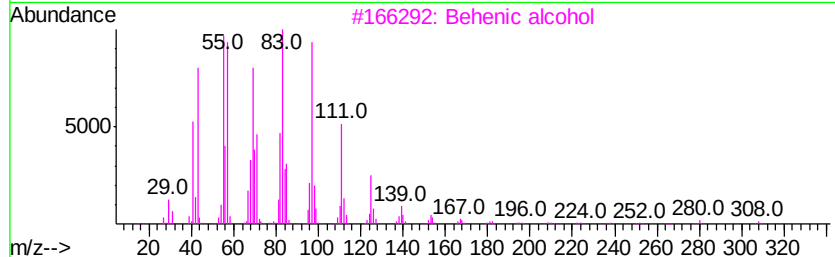
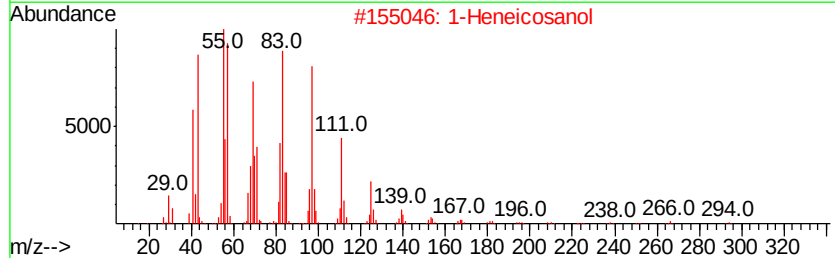
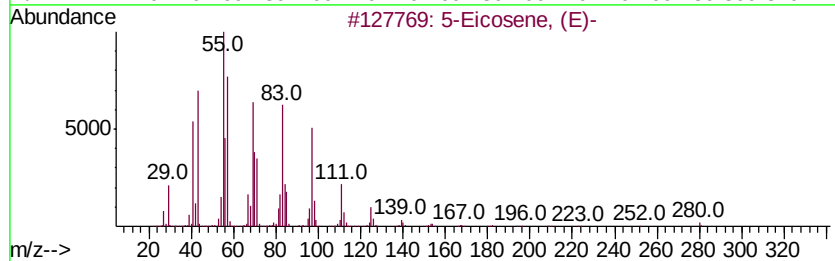
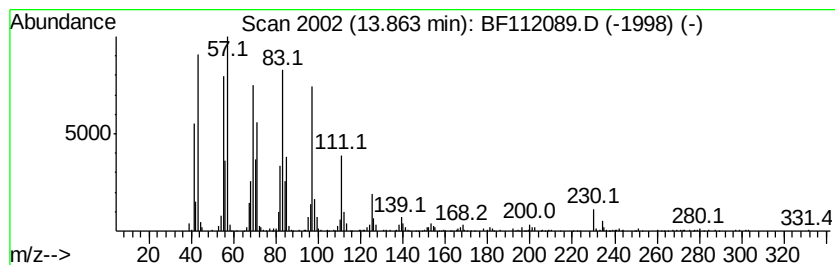
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 5-Eicosene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	6.88 ng	876512	Chrysene-d12	14.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2			1-Heneicosanol	312	C21H44O	015594-90-8	95
3			Behenic alcohol	326	C22H46O	000661-19-8	94
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94



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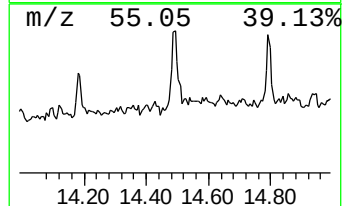
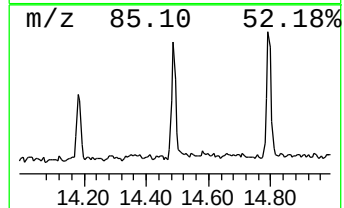
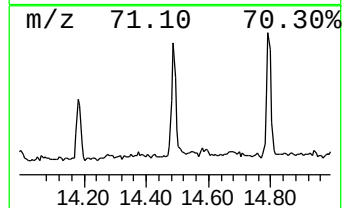
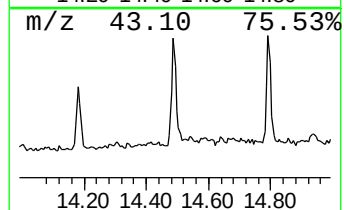
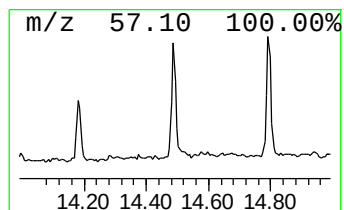
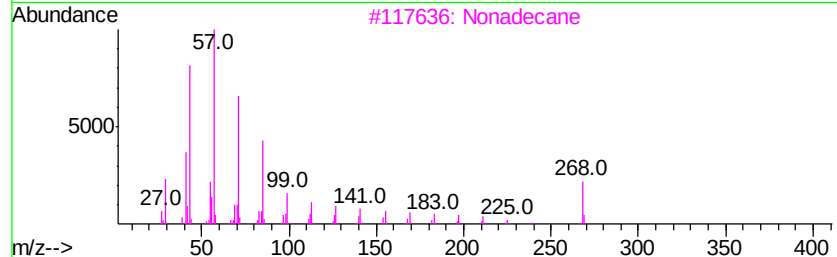
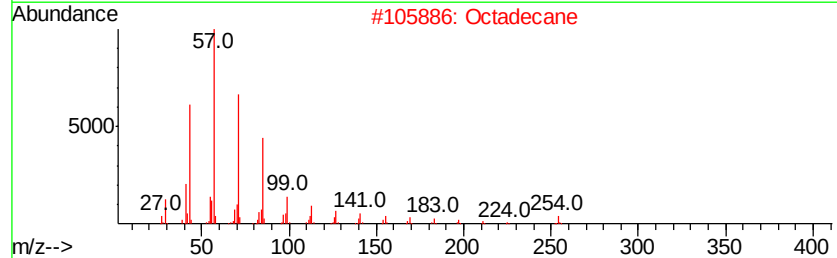
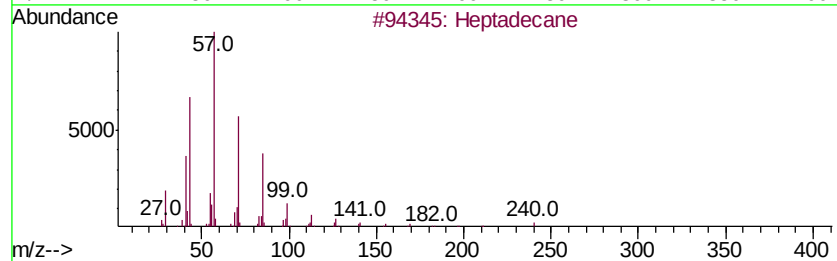
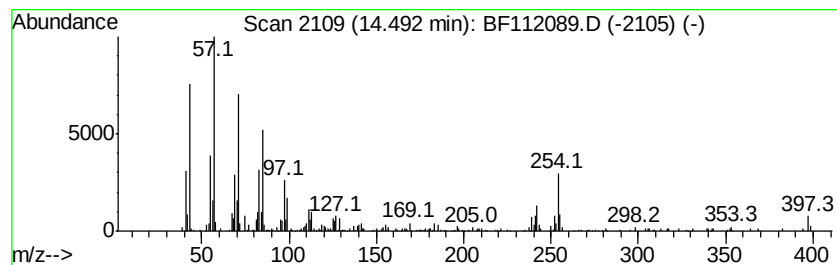
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.49	2.92 ng	372469	Chrysene-d12		14.03
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	96
2	Octadecane	254	C18H38	000593-45-3	91
3	Nonadecane	268	C19H40	000629-92-5	90
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
5	Tetracosane	338	C24H50	000646-31-1	87



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D-9

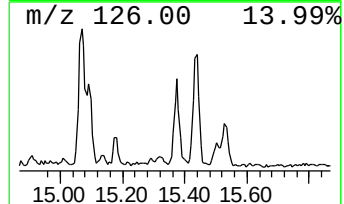
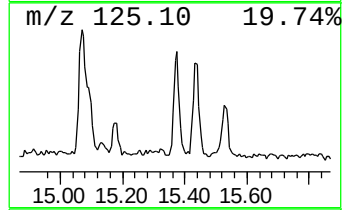
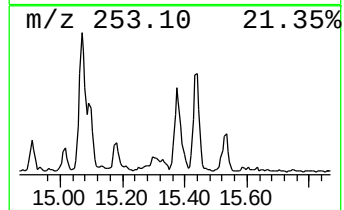
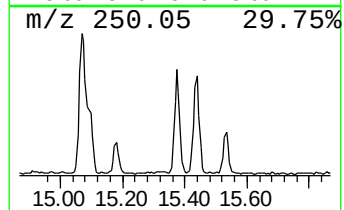
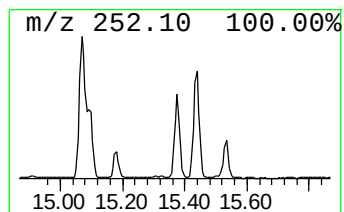
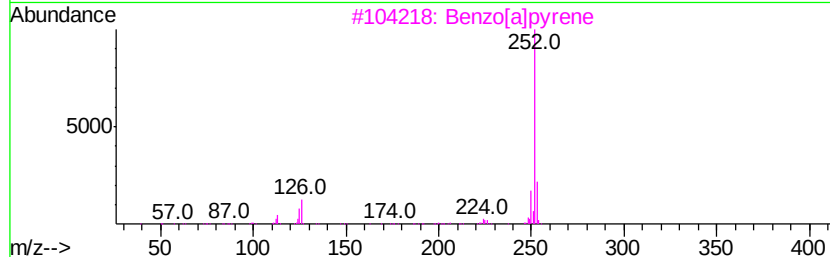
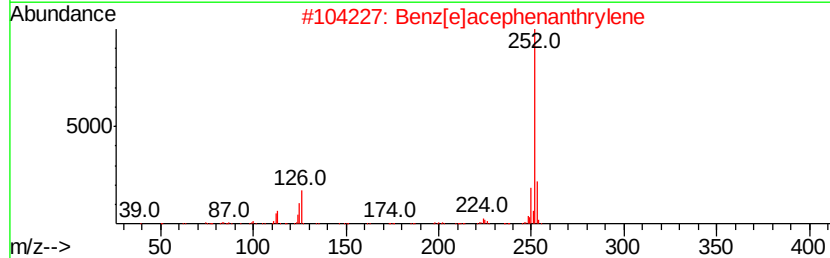
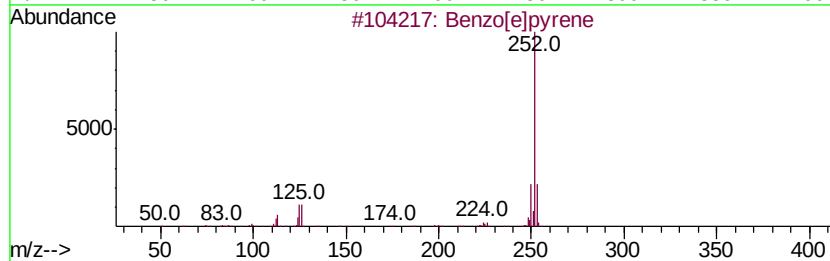
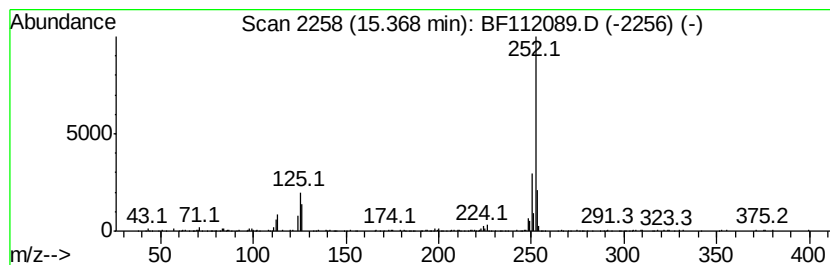
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	3.02 ng	344318	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]bpvrene	252	C20H12	000192-97-2	99
2		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	98
3		Benzo[a]bpvrene	252	C20H12	000050-32-8	94
4		Benzo[i]lfluoranthene	252	C20H12	000205-82-3	93
5		Perylene	252	C20H12	000198-55-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
Data File : BF112089.D
Acq On : 17 Jan 2019 15:11
Operator : JU/SJ
Sample : K1167-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
D-9

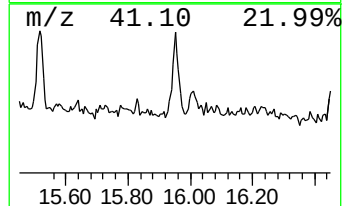
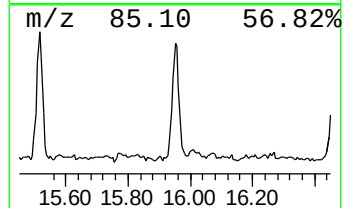
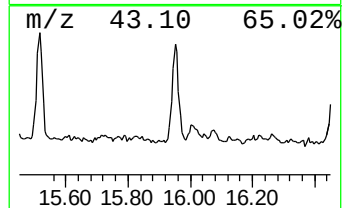
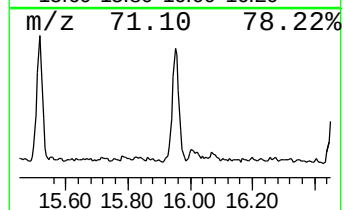
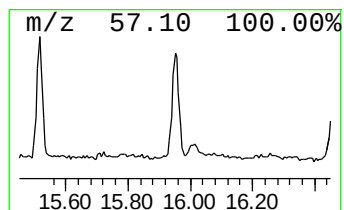
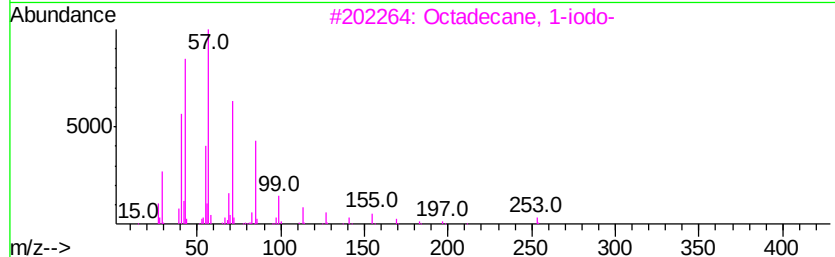
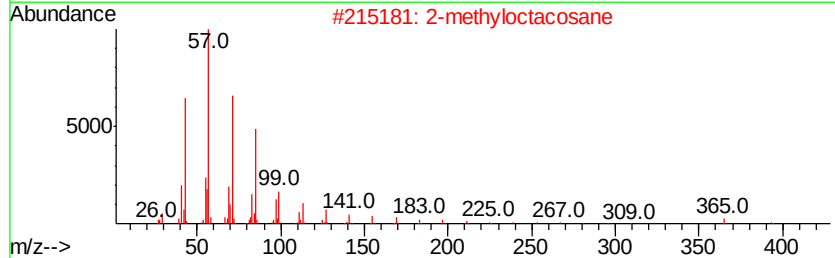
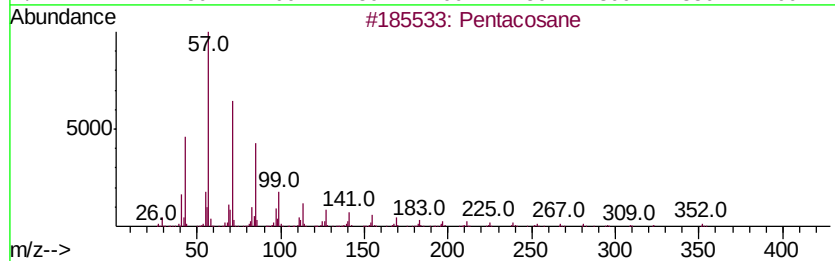
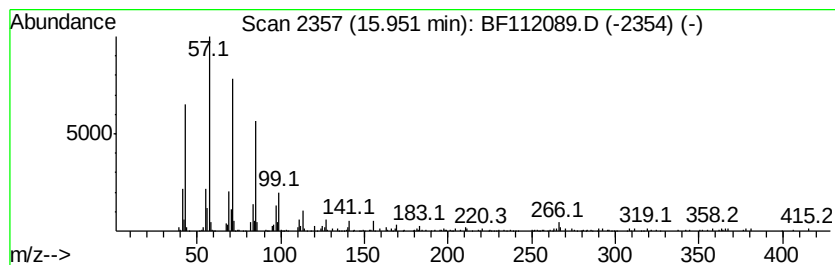
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pentacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.95	2.51 ng	286674	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacosane	352	C25H52	000629-99-2	89
2			2-methyloctacosane	408	C29H60	1000376-72-8	87
3			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86
4			2-Thiopheneacetic acid, 4-tetrad...	338	C20H34O2S	1000280-67-0	80
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF011719\
Data File : BF112089.D
Acq On : 17 Jan 2019 15:11
Operator : JU/SJ
Sample : K1167-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
D-9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF011619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butane, 2-methoxy...	2.15	2.8	ng	265742	1	6.86	1893480	20.0
2-Pentanone, 4-hy...	5.09	5.6	ng	527651	1	6.86	1893480	20.0
unknown6.64	6.64	88.8	ng	8404020	1	6.86	1893480	20.0
Benzene, 1-etheny...	6.70	2.9	ng	272531	1	6.86	1893480	20.0
Benzenemethanethi...	7.72	3.1	ng	367402	2	8.15	2387910	20.0
Benzene, 1-methyl...	8.36	2.4	ng	285828	2	8.15	2387910	20.0
o-Cymene	8.42	2.8	ng	331525	2	8.15	2387910	20.0
Anthracene, 1-met...	11.91	2.5	ng	297306	4	11.39	2364720	20.0
4H-Cyclopenta[def...	12.00	2.0	ng	239311	4	11.39	2364720	20.0
5-Eicosene, (E)-	13.86	6.9	ng	876512	5	14.03	2547280	20.0
Heptadecane	14.49	2.9	ng	372469	5	14.03	2547280	20.0
Benzo[e]pyrene	15.37	3.0	ng	344318	6	15.50	2281350	20.0
Pentacosane	15.95	2.5	ng	286674	6	15.50	2281350	20.0