

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030618\
 Data File : BF103476.D
 Acq On : 7 Mar 2018 00:34
 Operator : SJ/JU
 Sample : J1722-17
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-14

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:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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	2.116	3	5	13	rVB	198226	220797	6.31%	0.654%
2	5.104	511	513	527	rBV	47613	89025	2.54%	0.264%
3	5.498	576	580	607	rBV	2331378	3423991	97.87%	10.140%
4	6.510	748	752	771	rBV	2511744	3498489	100.00%	10.361%
5	6.645	771	775	796	rVB	2852834	3337078	95.39%	9.883%
6	6.869	810	813	830	rBV	726095	838091	23.96%	2.482%
7	7.028	836	840	855	rBV	2109704	2403050	68.69%	7.117%
8	7.439	906	910	934	rBV	1645240	2215777	63.34%	6.562%
9	8.157	1028	1032	1054	rVB	729107	1108809	31.69%	3.284%
10	9.227	1210	1214	1223	rBV	2742068	2871268	82.07%	8.503%
11	9.292	1223	1225	1229	rVB	71805	69290	1.98%	0.205%
12	9.622	1275	1281	1288	rBV	160614	228836	6.54%	0.678%
13	9.774	1303	1307	1311	rBV	39410	54847	1.57%	0.162%
14	9.822	1312	1315	1319	rVB	150370	128189	3.66%	0.380%
15	9.910	1326	1330	1334	rBV2	1001435	999739	28.58%	2.961%
16	10.127	1361	1367	1368	rVV2	32362	50044	1.43%	0.148%
17	10.145	1368	1370	1374	rVB3	45424	47724	1.36%	0.141%
18	10.321	1397	1400	1403	rVB	193624	145590	4.16%	0.431%
19	10.469	1422	1425	1430	rVB	43609	52370	1.50%	0.155%
20	10.545	1433	1438	1440	rBV	51299	70022	2.00%	0.207%
21	10.710	1461	1466	1478	rBV	1195861	1547665	44.24%	4.583%
22	10.798	1478	1481	1490	rVB	144965	192696	5.51%	0.571%
23	10.945	1503	1506	1511	rBV	40124	46677	1.33%	0.138%
24	11.245	1555	1557	1560	rBV2	82888	65068	1.86%	0.193%
25	11.404	1580	1584	1587	rBV	812724	918993	26.27%	2.722%
26	11.427	1587	1588	1595	rVV	565965	526779	15.06%	1.560%
27	11.486	1595	1598	1602	rVV	125855	131845	3.77%	0.390%
28	11.651	1622	1626	1628	rBV2	56376	66747	1.91%	0.198%
29	11.910	1667	1670	1673	rBV2	112493	144671	4.14%	0.428%
30	11.939	1673	1675	1680	rVB2	103011	107573	3.07%	0.319%
31	11.986	1681	1683	1686	rVB	48315	41356	1.18%	0.122%
32	12.021	1686	1689	1692	rBV	113325	142730	4.08%	0.423%
33	12.204	1717	1720	1723	rBV	61055	61272	1.75%	0.181%
34	12.245	1725	1727	1731	rVB2	45321	42283	1.21%	0.125%

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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

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35	12.457	1760	1763	1768	rBV3	54415	85275	2.44%	0.253%
36	12.627	1788	1792	1800	rBV	930692	912359	26.08%	2.702%
37	12.792	1815	1820	1824	rBV5	56340	108229	3.09%	0.321%
38	12.857	1824	1831	1836	rBV	772705	959766	27.43%	2.842%
39	12.992	1849	1854	1857	rBV	2017872	1953839	55.85%	5.786%
40	13.080	1864	1869	1873	rBV2	53063	95653	2.73%	0.283%
41	13.162	1880	1883	1884	rBV2	52469	50591	1.45%	0.150%
42	13.186	1885	1887	1892	rVV2	107890	122389	3.50%	0.362%
43	13.251	1896	1898	1901	rVV	71809	62968	1.80%	0.186%
44	13.292	1903	1905	1912	rVB2	71290	99308	2.84%	0.294%
45	13.421	1924	1927	1929	rBV2	67376	79075	2.26%	0.234%
46	13.633	1961	1963	1965	rBV2	43297	38129	1.09%	0.113%
47	13.704	1973	1975	1978	rVB	81612	73414	2.10%	0.217%
48	13.810	1991	1993	1995	rBV	70335	60454	1.73%	0.179%
49	13.868	2000	2003	2009	rBV4	361256	475179	13.58%	1.407%
50	14.062	2031	2036	2039	rBV2	740623	1092773	31.24%	3.236%
51	14.086	2039	2040	2047	rVB	355910	307026	8.78%	0.909%
52	14.168	2052	2054	2056	rBV	62355	53883	1.54%	0.160%
53	14.557	2118	2120	2123	rBV3	73036	72274	2.07%	0.214%
54	15.133	2214	2218	2221	rBV	254171	438164	12.52%	1.298%
55	15.445	2268	2271	2275	rBV	172014	187037	5.35%	0.554%
56	15.509	2279	2282	2286	rVV	202845	242233	6.92%	0.717%
57	15.574	2290	2293	2297	rBV	256807	307833	8.80%	0.912%

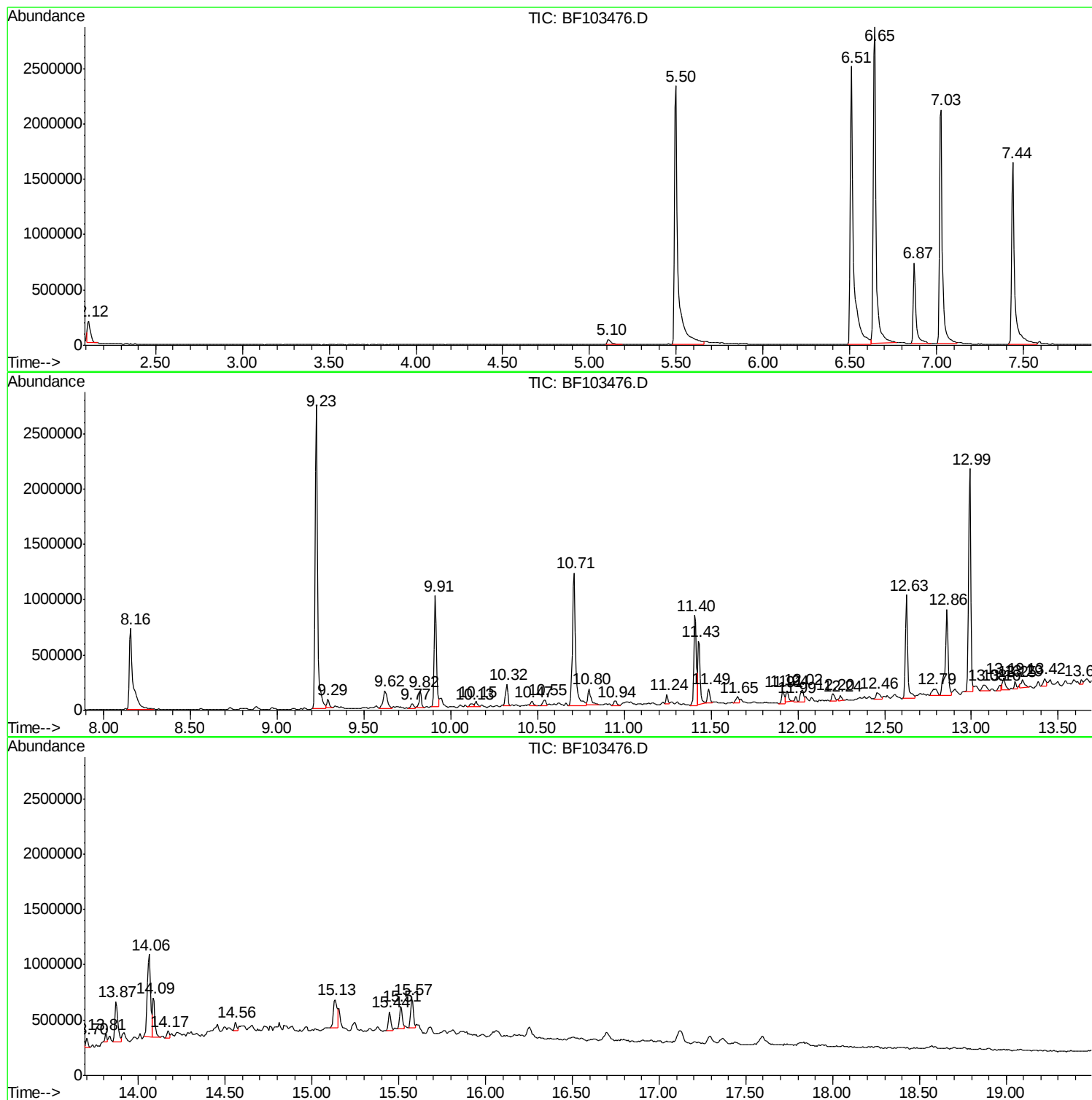
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Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : J1722-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-14

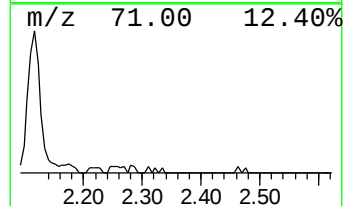
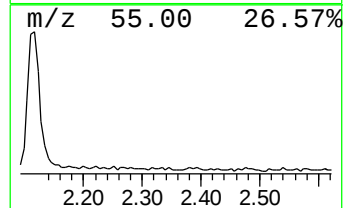
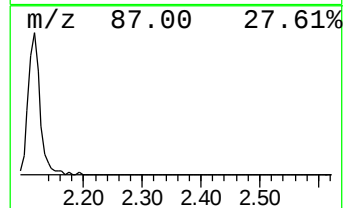
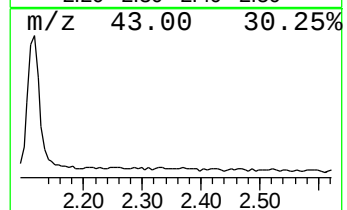
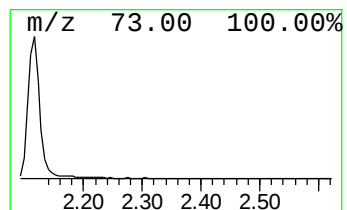
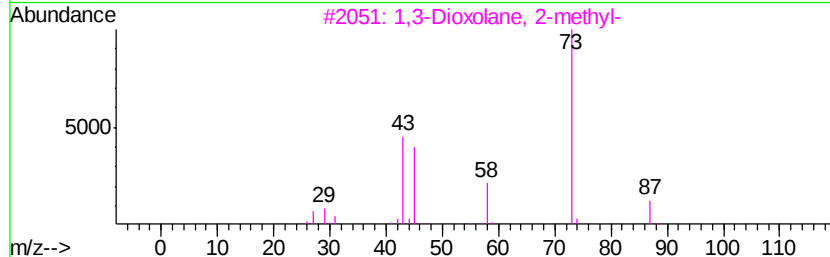
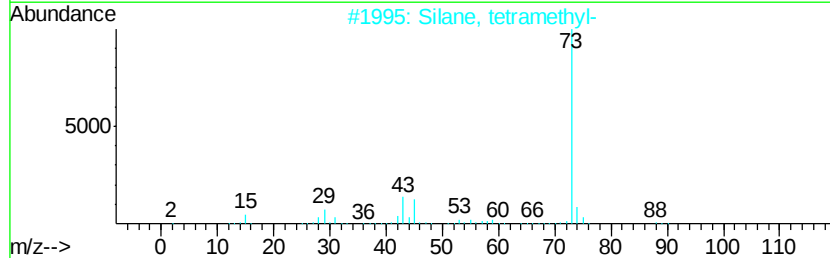
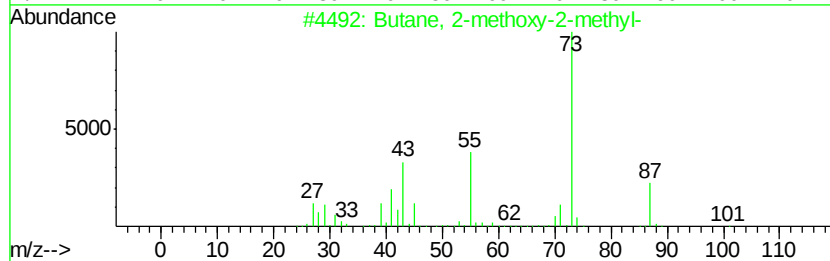
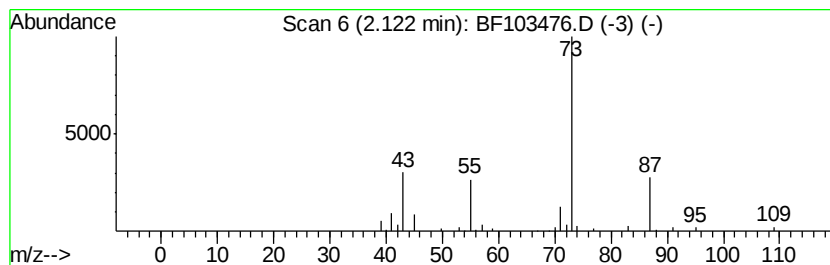
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	5.27 ng	220797	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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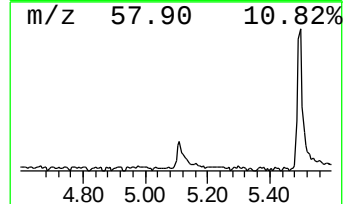
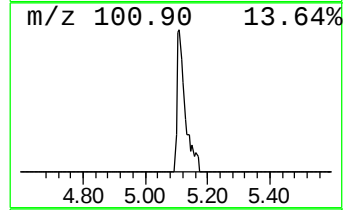
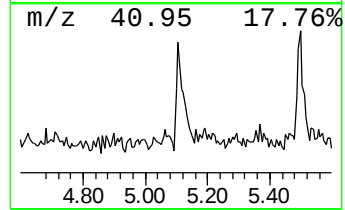
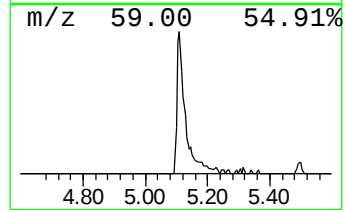
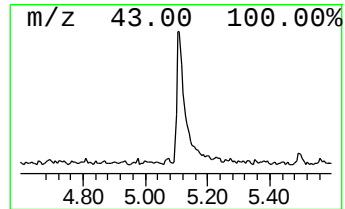
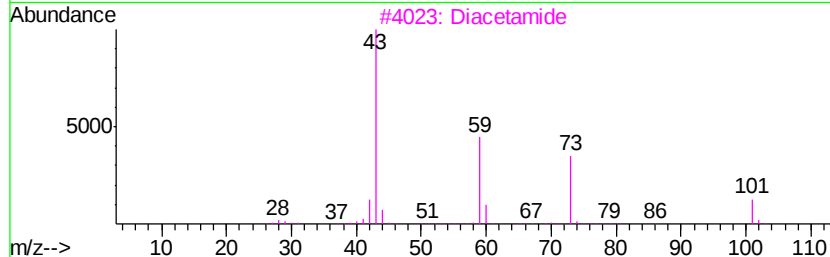
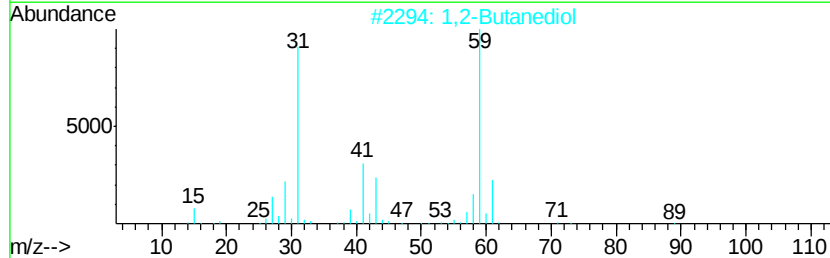
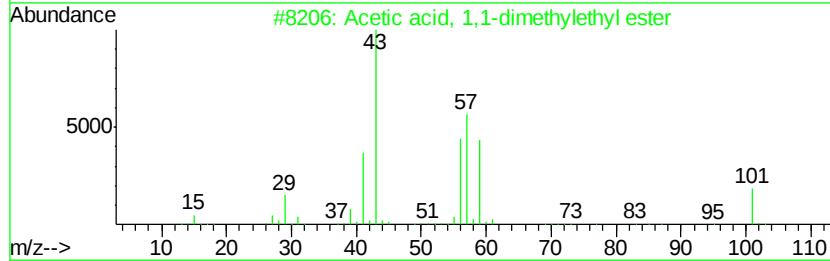
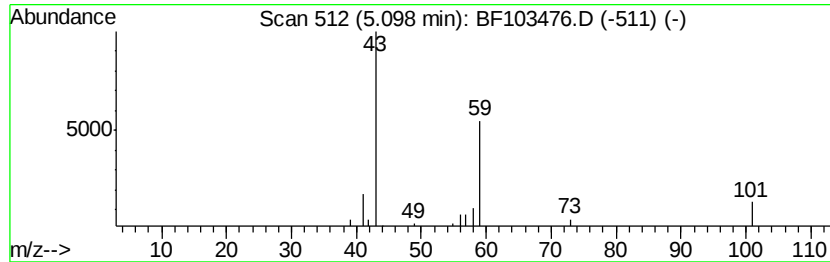
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.10 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	2.12 ng	89025	1,4-Dichlorobenzene-d4	6.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
2			1,2-Butanediol	90	C4H10O2	000584-03-2	37
3			Diacetamide	101	C4H7NO2	000625-77-4	36
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25
5			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25



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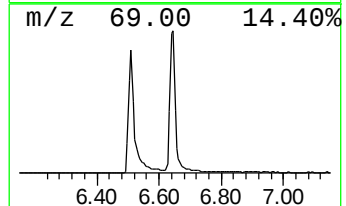
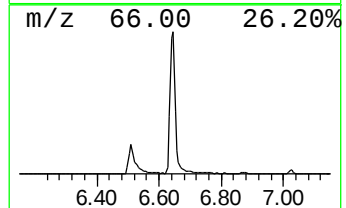
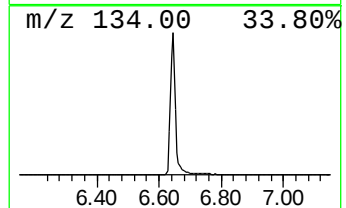
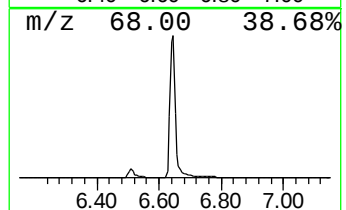
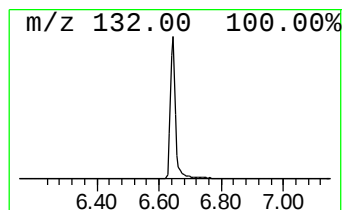
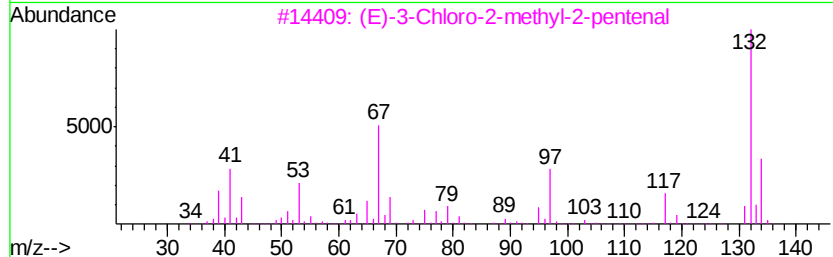
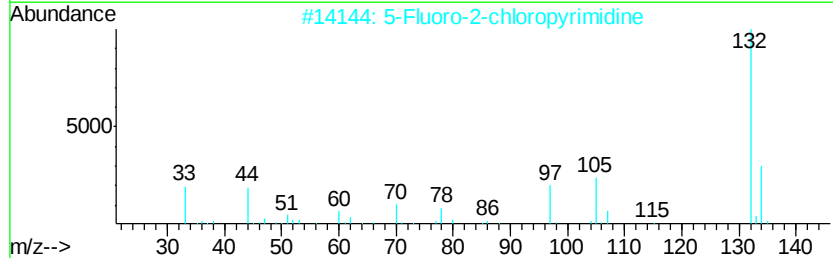
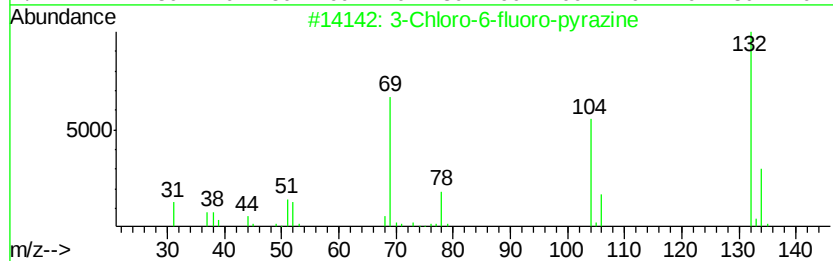
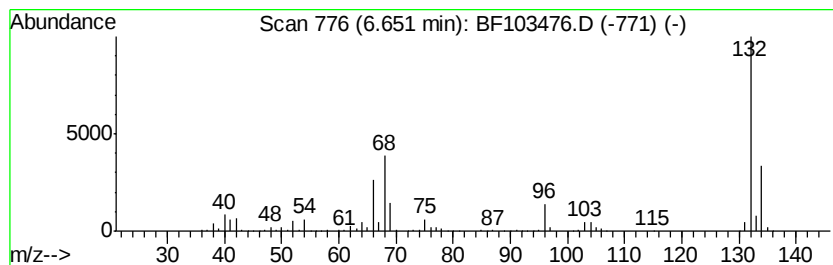
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	79.64 ng	3337080	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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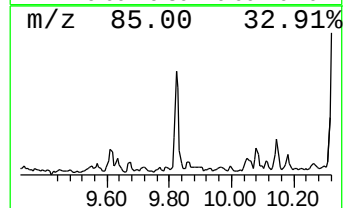
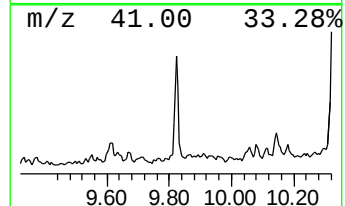
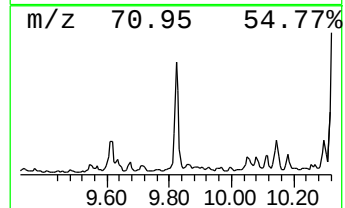
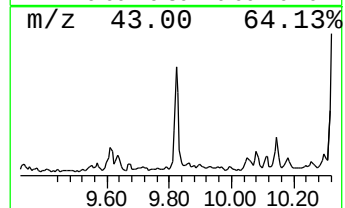
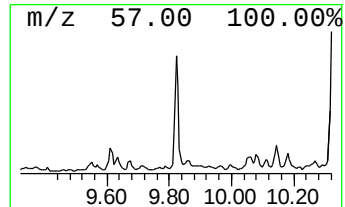
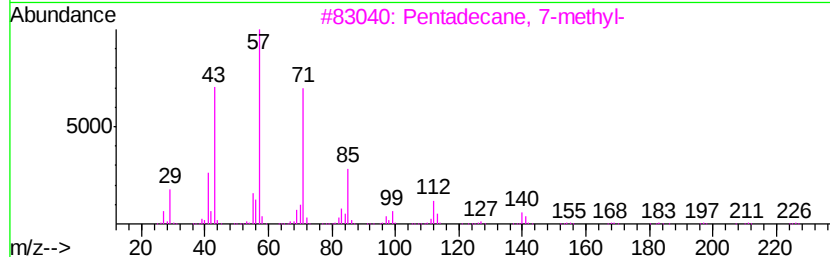
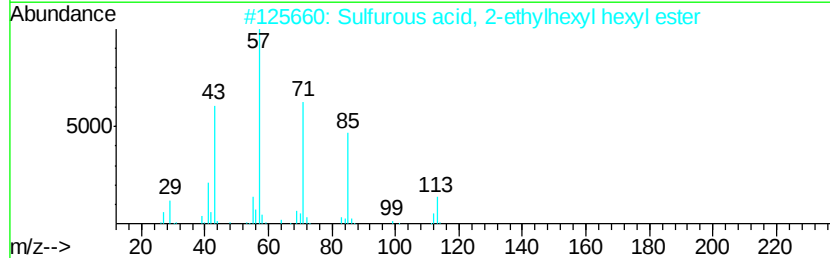
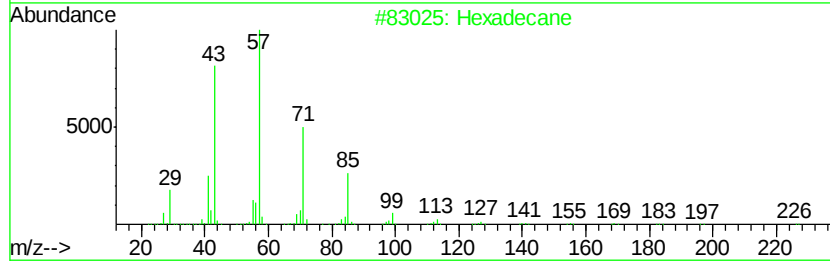
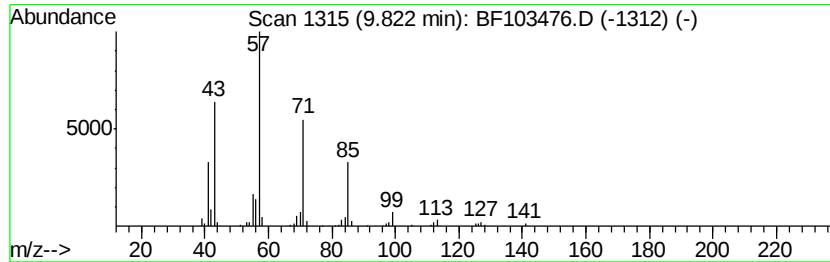
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	2.56 ng	128189	Acenaphthene-d10	9.91

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	87
2	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	80
3	Pentadecane, 7-methyl-		226	C16H34	006165-40-8	78
4	Octane, 3-ethyl-2,7-dimethyl-		170	C12H26	062183-55-5	59
5	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	59



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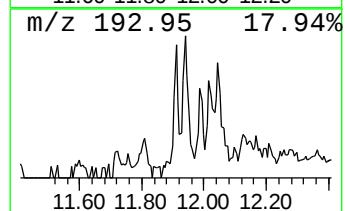
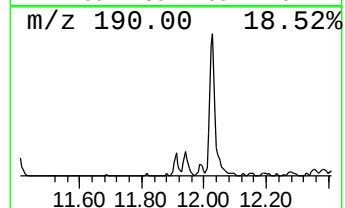
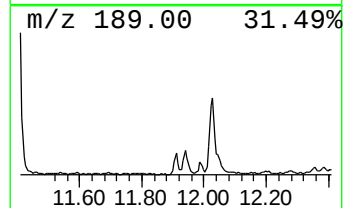
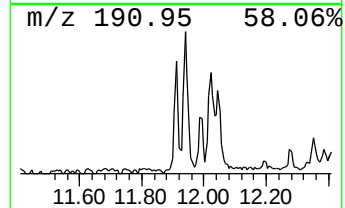
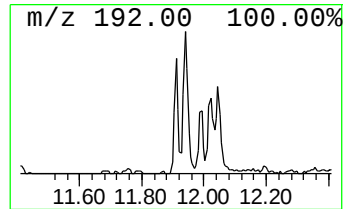
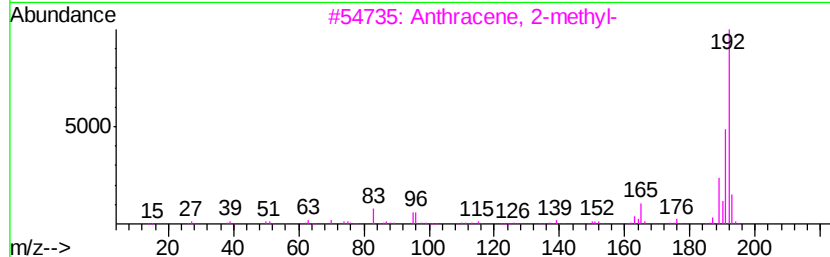
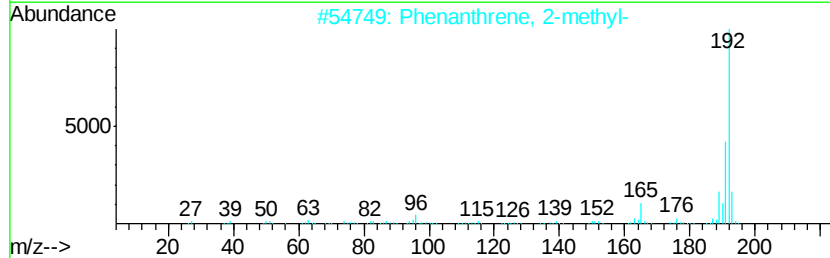
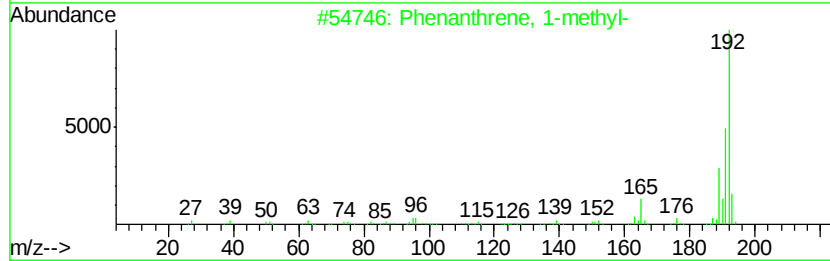
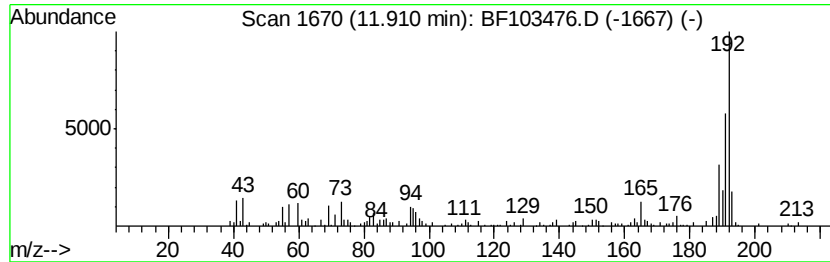
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.91	3.15 ng	144671	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
4	1H-Cyclopropa[1,1]phenanthrene,1a,...	192	C15H12	000949-41-7	86
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	76



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103476.D
Acq On : 7 Mar 2018 00:34
Operator : SJ/JU
Sample : J1722-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SP-14

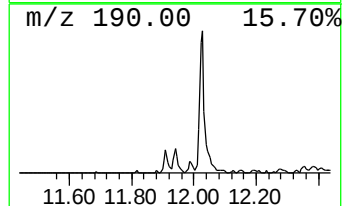
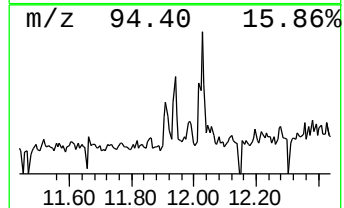
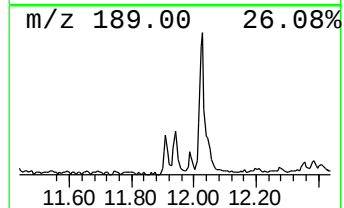
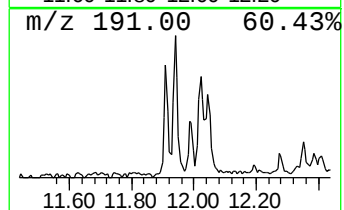
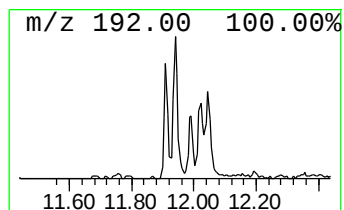
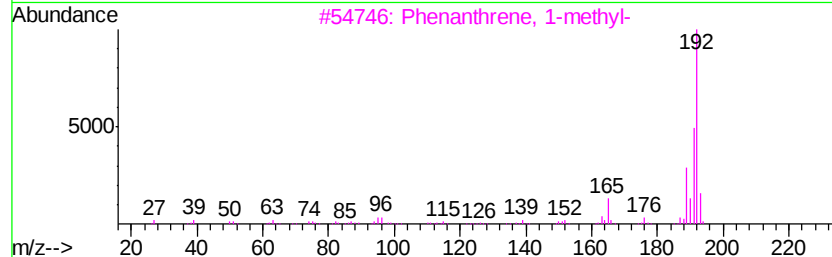
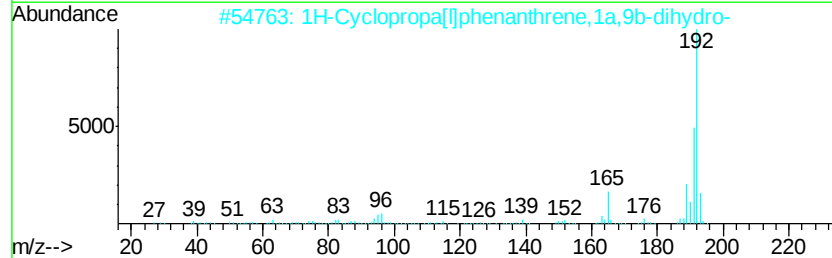
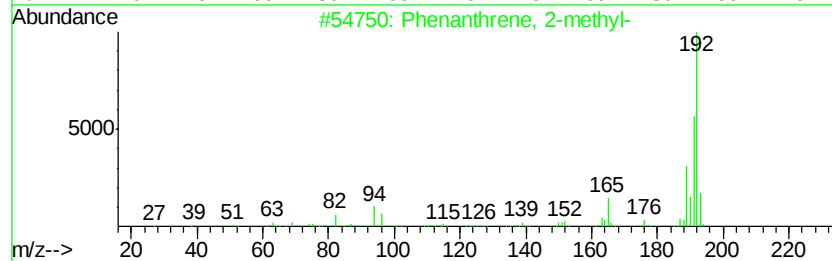
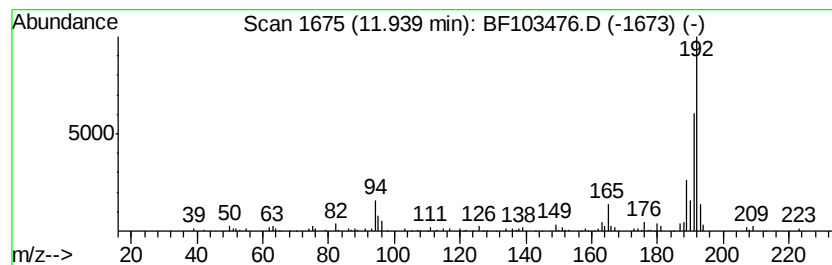
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	2.34 ng	107573	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103476.D
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Misc :
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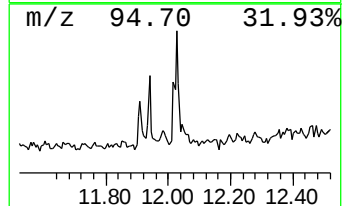
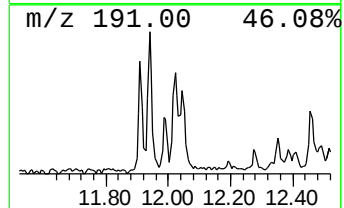
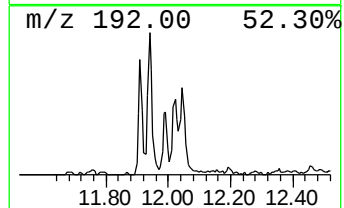
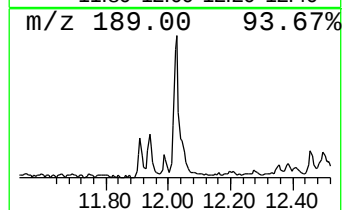
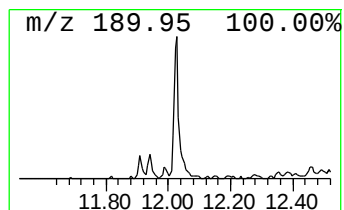
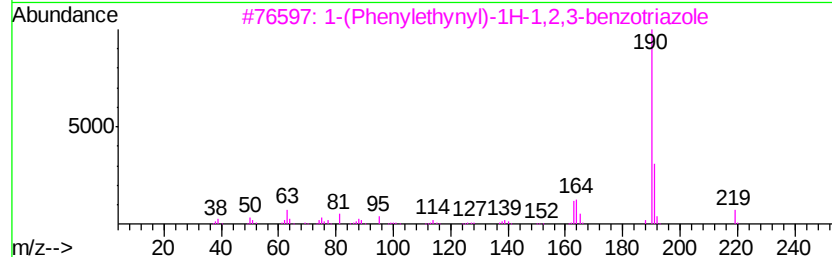
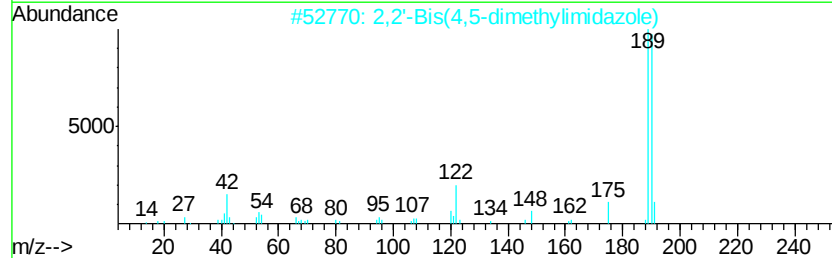
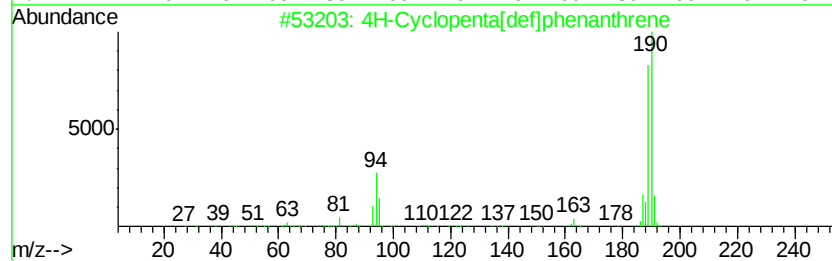
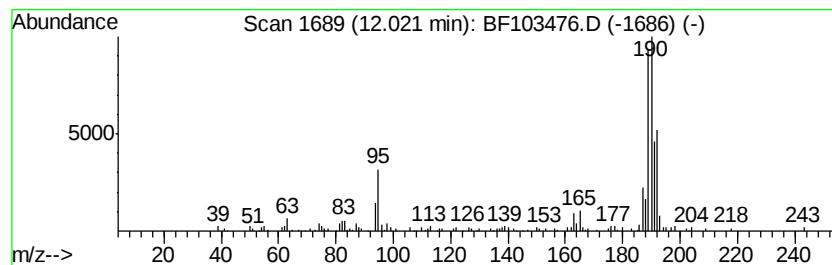
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	3.11 ng	142730	Phenanthrene-d10	11.40

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
2			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
3			1-(Phenylethynyl)-1H-1,2,3-benzo...	219	C14H9N3	209912-23-2	25
4			Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	25
5			Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	18



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103476.D
Acq On : 7 Mar 2018 00:34
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Misc :
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Instrument :
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ClientSampleId :
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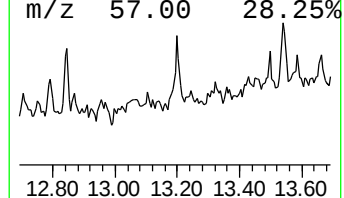
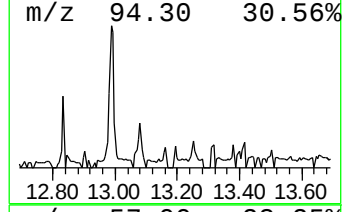
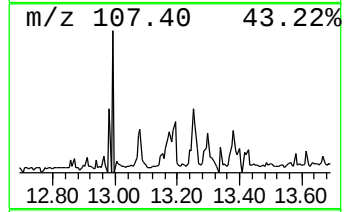
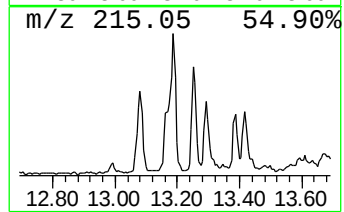
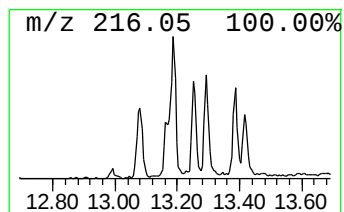
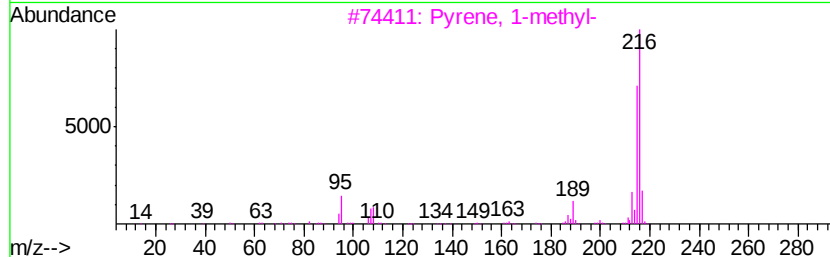
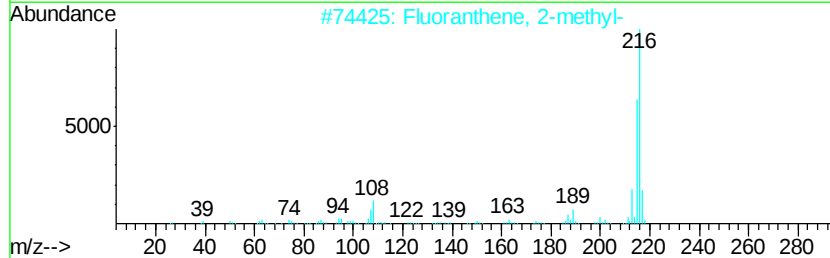
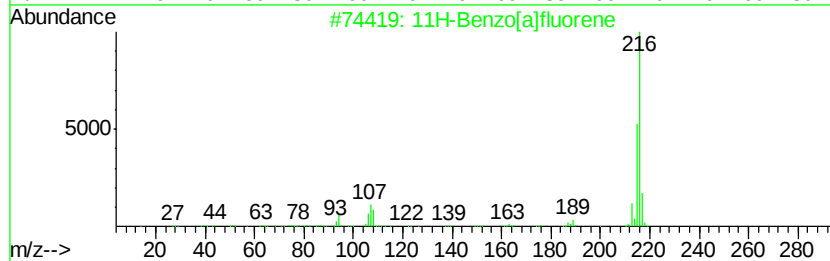
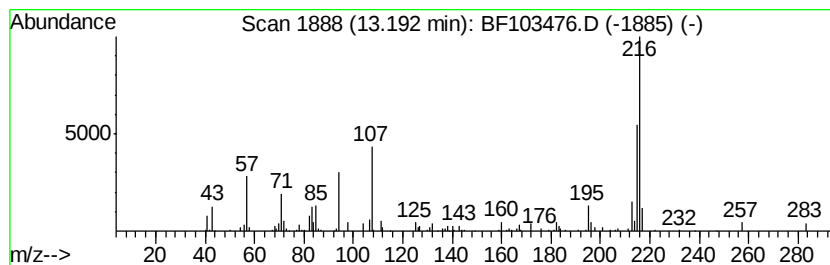
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[a]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	2.24 ng	122389	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
2		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	53
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	53
4		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	53
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	50



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103476.D
Acq On : 7 Mar 2018 00:34
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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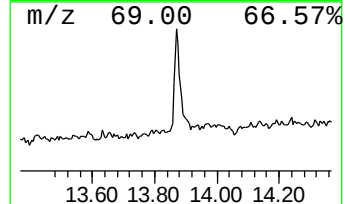
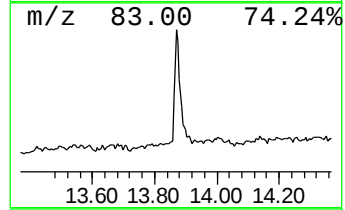
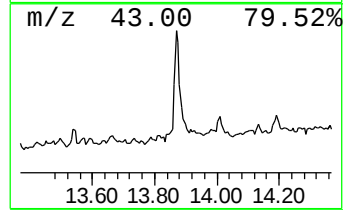
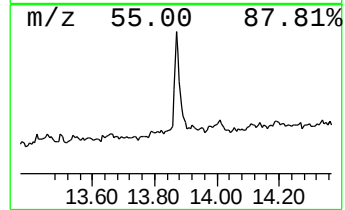
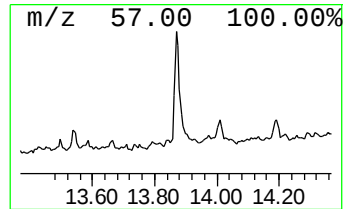
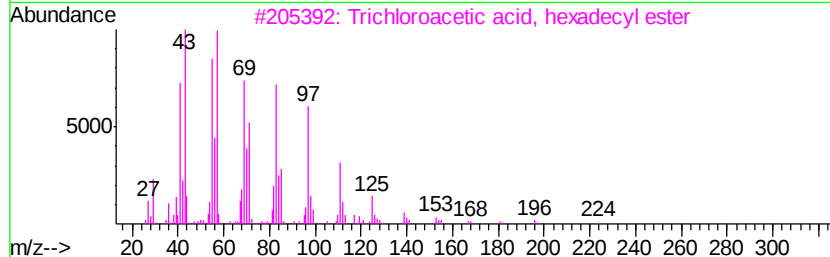
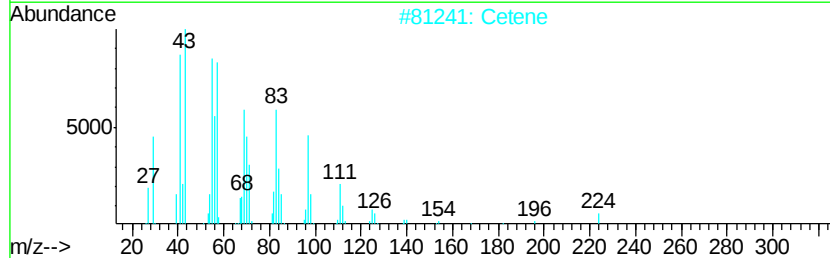
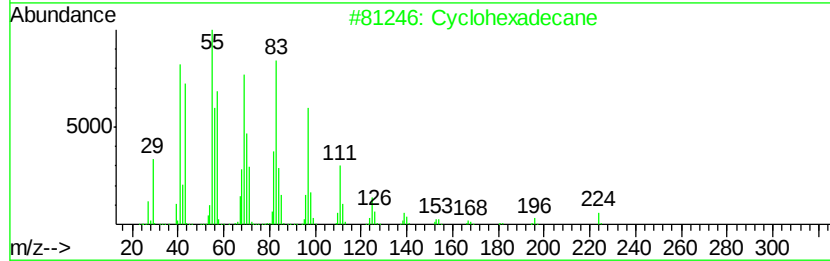
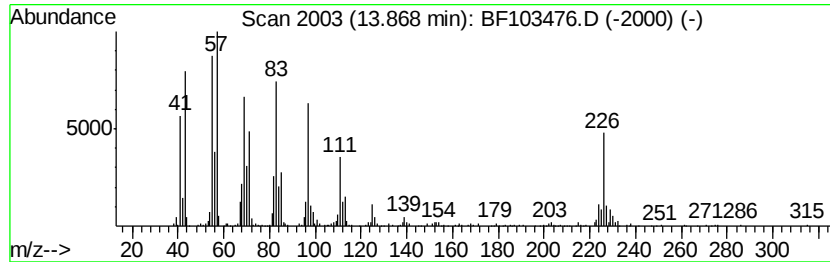
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cyclohexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	8.70 ng	475179	Chrysene-d12	14.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	92
2			Cetene	224	C16H32	000629-73-2	86
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	76
4			1-Heptacosanol	396	C27H56O	002004-39-9	76
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	76



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103476.D
Acq On : 7 Mar 2018 00:34
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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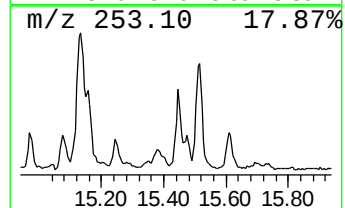
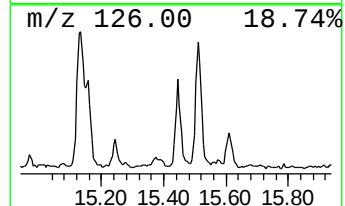
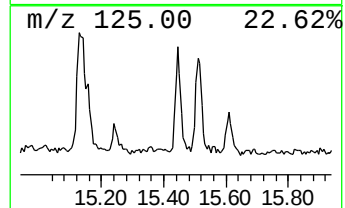
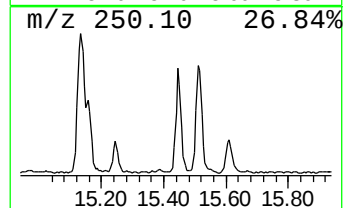
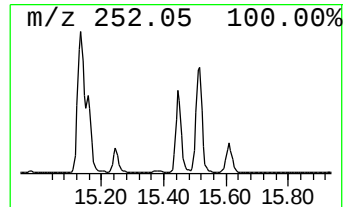
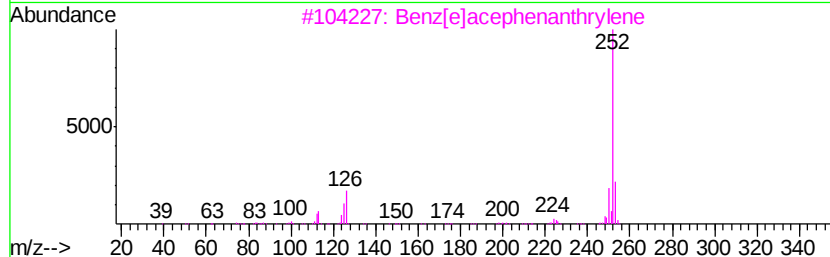
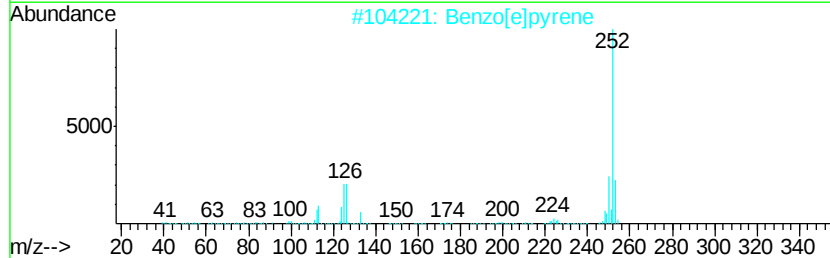
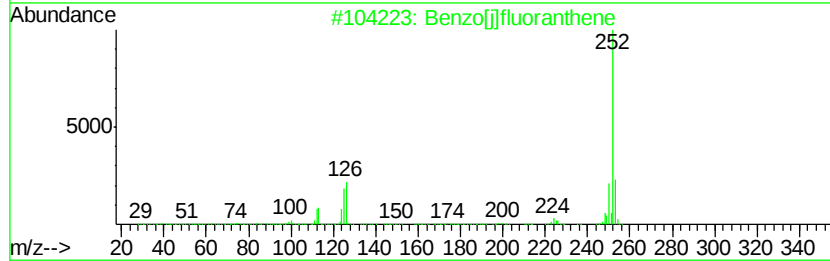
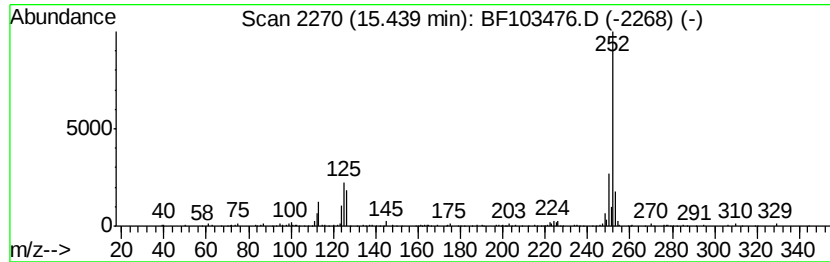
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzo[j]fluoranthene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.44	12.15 ng	187037	Perylene-d12	15.57

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



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030618\
Data File : BF103476.D
Acq On : 7 Mar 2018 00:34
Operator : SJ/JU
Sample : J1722-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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
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unknown5.10	5.10	2.1	ng	89025	1	6.87	838091	20.0
unknown6.65	6.65	79.6	ng	3337080	1	6.87	838091	20.0
Hexadecane	9.82	2.6	ng	128189	3	9.91	999739	20.0
Phenanthrene, 1-m...	11.91	3.1	ng	144671	4	11.40	918993	20.0
Phenanthrene, 2-m...	11.94	2.3	ng	107573	4	11.40	918993	20.0
4H-Cyclopenta[def...	12.02	3.1	ng	142730	4	11.40	918993	20.0
11H-Benzo[a]fluorene	13.19	2.2	ng	122389	5	14.06	1092770	20.0
Cyclohexadecane	13.87	8.7	ng	475179	5	14.06	1092770	20.0
Benzo[j]fluoranthene	15.44	12.2	ng	187037	6	15.57	307833	20.0