

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
 Data File : BF098969.D
 Acq On : 30 Sep 2017 18:15
 Operator : SJ/JU
 Sample : I5563-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYSEG-PH4-SB-ESMI-6

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-BF092317.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	2.234	20	25	33	rVB	180735	245872	6.69%	0.569%
2	5.151	517	521	536	rVB	404851	590765	16.07%	1.367%
3	5.545	580	588	591	rBV	3412485	3512422	95.55%	8.127%
4	6.545	752	758	765	rBV	2897497	3550809	96.59%	8.216%
5	6.687	777	782	786	rBV	3331406	3546963	96.49%	8.207%
6	6.916	817	821	827	rBV	1326894	1061944	28.89%	2.457%
7	7.075	842	848	851	rBV	3170862	2884744	78.48%	6.675%
8	7.481	911	917	920	rBV	2266552	2233000	60.75%	5.167%
9	8.198	1035	1039	1041	rBV	1647288	1343804	36.56%	3.109%
10	8.216	1041	1042	1045	rVB	206777	149012	4.05%	0.345%
11	8.910	1156	1160	1165	rVB	63127	51946	1.41%	0.120%
12	9.010	1173	1177	1181	rBV	46765	43194	1.18%	0.100%
13	9.275	1216	1222	1225	rBV	3888012	3675998	100.00%	8.506%
14	9.369	1236	1238	1243	rVB2	39973	38497	1.05%	0.089%
15	9.610	1275	1279	1281	rBV	40560	43040	1.17%	0.100%
16	9.663	1285	1288	1295	rVB	586458	483521	13.15%	1.119%
17	9.816	1311	1314	1317	rVB	72806	60088	1.63%	0.139%
18	9.875	1319	1324	1327	rVB2	40210	40817	1.11%	0.094%
19	9.957	1331	1338	1341	rVV	1316015	1305313	35.51%	3.020%
20	9.986	1341	1343	1346	rVB	177368	129683	3.53%	0.300%
21	10.157	1366	1372	1375	rBV	100027	106725	2.90%	0.247%
22	10.375	1402	1409	1412	rBV2	90324	115800	3.15%	0.268%
23	10.498	1427	1430	1435	rVB	152811	143707	3.91%	0.333%
24	10.657	1451	1457	1464	rVB3	39852	58762	1.60%	0.136%
25	10.745	1467	1472	1475	rBV	1526753	1448033	39.39%	3.351%
26	10.845	1486	1489	1498	rBV	140382	174349	4.74%	0.403%
27	11.045	1518	1523	1526	rBV5	48690	77984	2.12%	0.180%
28	11.245	1554	1557	1560	rBV	48729	40270	1.10%	0.093%
29	11.292	1562	1565	1568	rVV2	102604	99454	2.71%	0.230%
30	11.339	1570	1573	1577	rVB	81143	83796	2.28%	0.194%
31	11.445	1587	1591	1593	rBV	1097181	1108658	30.16%	2.565%
32	11.469	1593	1595	1598	rVV	1338535	1003656	27.30%	2.322%
33	11.516	1600	1603	1606	rVV	343756	276430	7.52%	0.640%
34	11.551	1606	1609	1614	rVB	65212	72967	1.98%	0.169%

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35	11.669	1623	1629	1635	rBV	162901	208213	5.66%	0.482%
36	11.945	1672	1676	1677	rBV	113974	111475	3.03%	0.258%
37	11.969	1677	1680	1684	rVB	184973	182737	4.97%	0.423%
38	12.016	1685	1688	1691	rBV	67927	60984	1.66%	0.141%
39	12.057	1691	1695	1697	rBV2	236441	239327	6.51%	0.554%
40	12.074	1697	1698	1702	rVB	85135	64040	1.74%	0.148%
41	12.127	1704	1707	1709	rBV2	39221	39742	1.08%	0.092%
42	12.239	1723	1726	1728	rBV	107827	98675	2.68%	0.228%
43	12.263	1728	1730	1733	rVB	137695	107664	2.93%	0.249%
44	12.492	1767	1769	1772	rBV	55866	60826	1.65%	0.141%
45	12.580	1782	1784	1788	rVB	69751	65425	1.78%	0.151%
46	12.663	1794	1798	1801	rBV	1845625	1631134	44.37%	3.774%
47	12.886	1830	1836	1841	rBV	1512200	1708060	46.47%	3.952%
48	12.933	1842	1844	1849	rVB2	63430	69665	1.90%	0.161%
49	13.027	1853	1860	1863	rBV	2451693	2333535	63.48%	5.400%
50	13.210	1889	1891	1895	rVB	196045	183821	5.00%	0.425%
51	13.280	1900	1903	1905	rBV	158119	134975	3.67%	0.312%
52	13.316	1907	1909	1912	rVB2	108275	90890	2.47%	0.210%
53	13.715	1975	1977	1983	rVB	92407	102561	2.79%	0.237%
54	13.833	1994	1997	1999	rBV3	110228	111873	3.04%	0.259%
55	13.857	1999	2001	2004	rBV	88150	69158	1.88%	0.160%
56	13.886	2004	2006	2007	rBV	89414	73671	2.00%	0.170%
57	13.904	2007	2009	2012	rVV3	107300	108854	2.96%	0.252%
58	14.080	2035	2039	2042	rBV2	1130775	1620723	44.09%	3.750%
59	14.110	2042	2044	2050	rVB	790418	694788	18.90%	1.608%
60	14.192	2056	2058	2061	rVB	130985	101315	2.76%	0.234%
61	15.139	2215	2219	2223	rBV	600496	1014223	27.59%	2.347%
62	15.457	2269	2273	2278	rVB	333898	444797	12.10%	1.029%
63	15.521	2280	2284	2289	rVB	488906	602490	16.39%	1.394%
64	15.586	2290	2295	2298	rBV	581856	743958	20.24%	1.721%
65	17.092	2548	2551	2558	rVB3	186391	335254	9.12%	0.776%

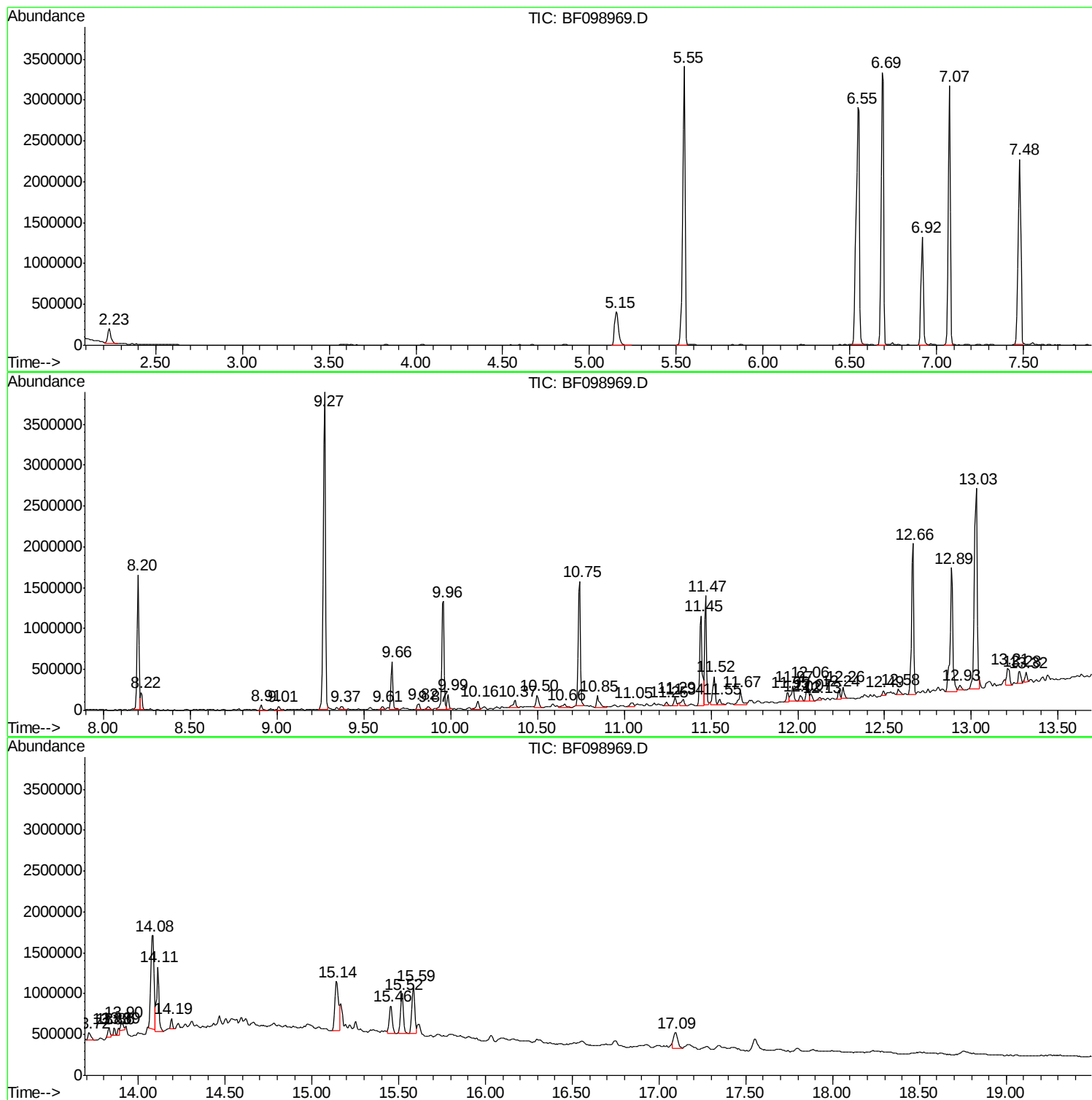
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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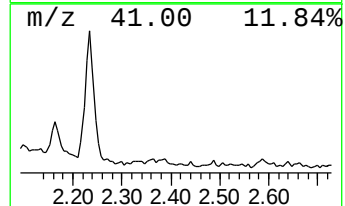
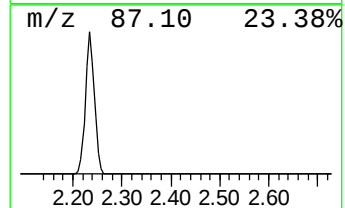
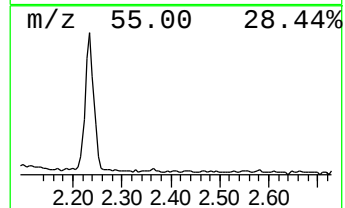
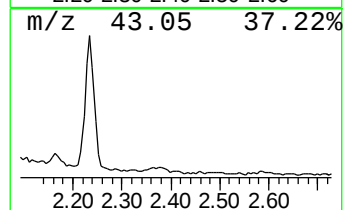
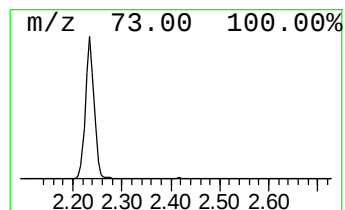
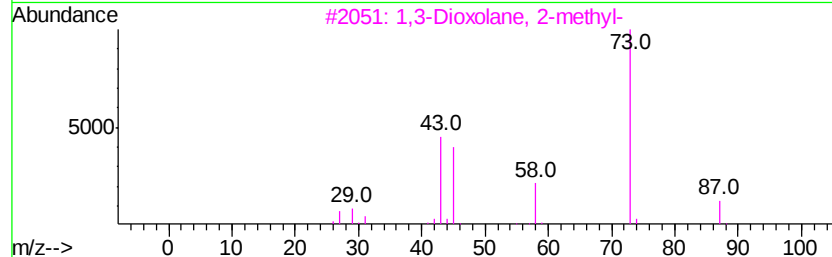
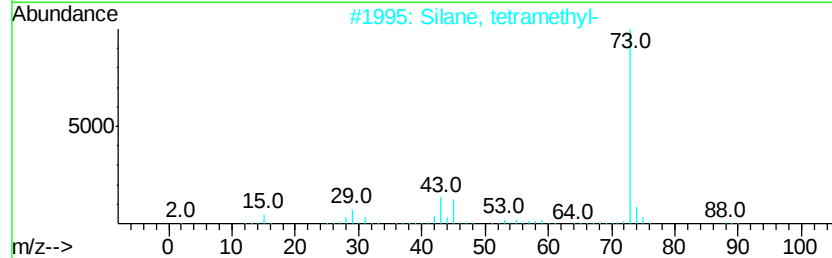
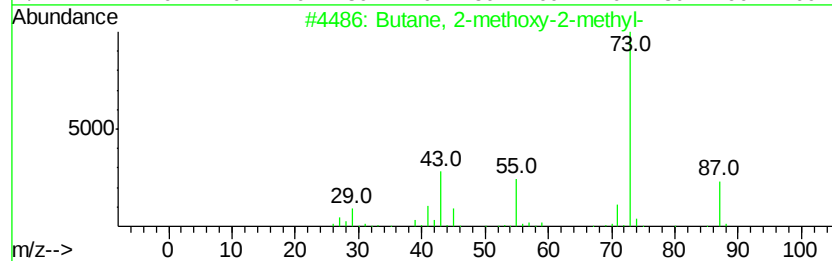
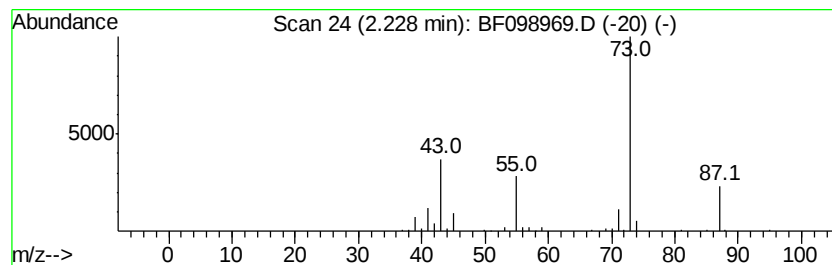
Instrument :
BNA_F
ClientSampleId :
NYSEG-PH4-SB-ESMI-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.23	4.63 ng	245872	1,4-Dichlorobenzene-d4	6.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4	3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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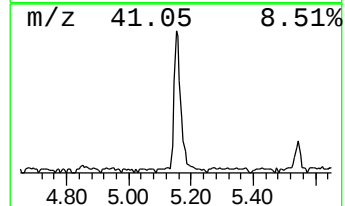
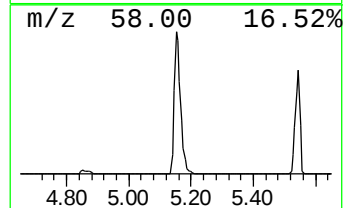
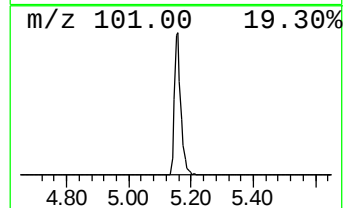
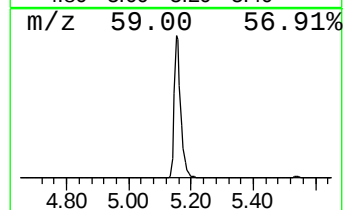
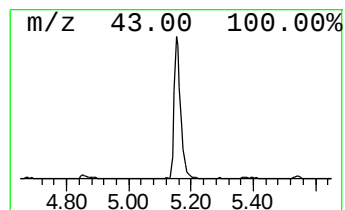
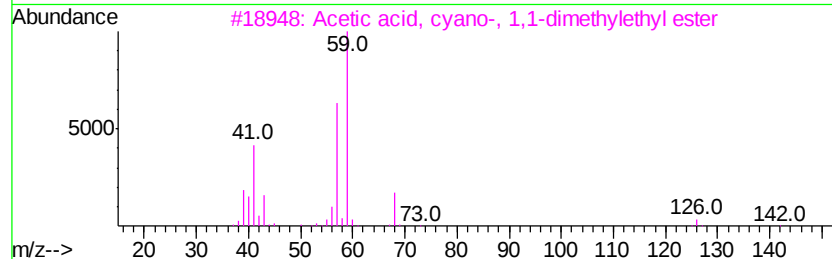
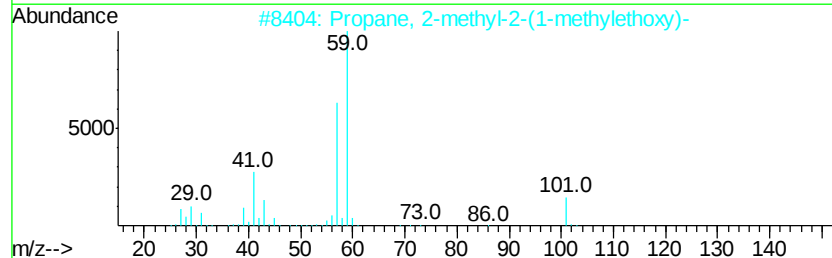
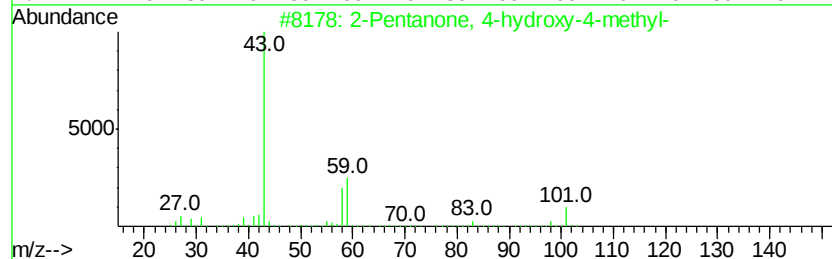
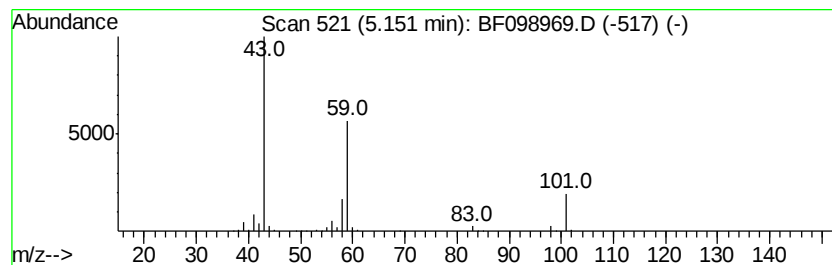
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.15	11.13 ng	590765	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	36
3		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		2-Butanone, 3-ethoxy-3-methyl-	130	C7H14O2	036687-99-7	9



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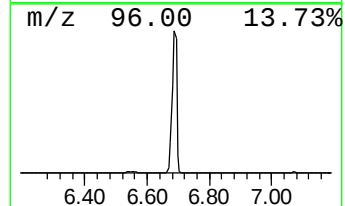
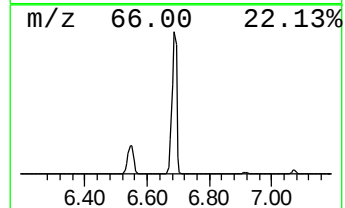
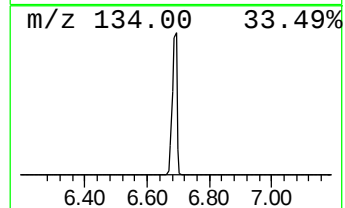
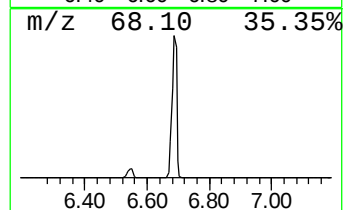
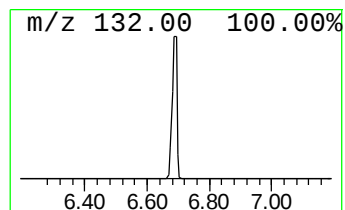
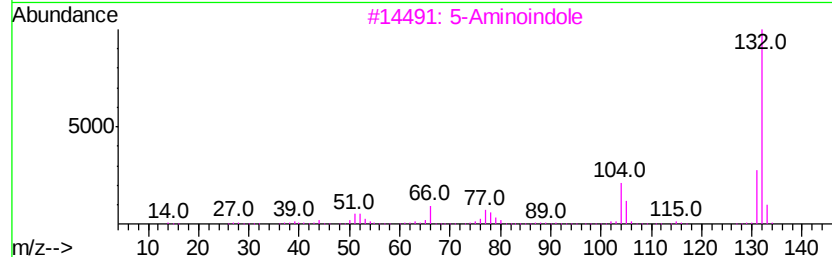
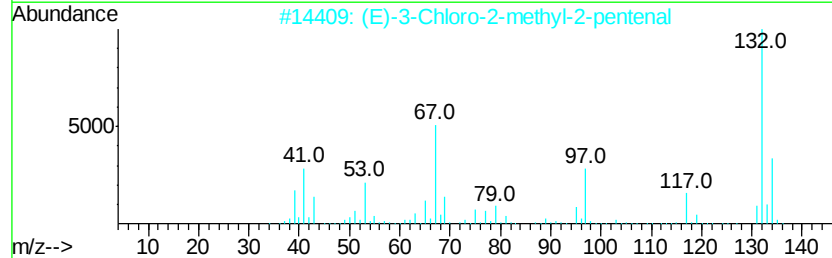
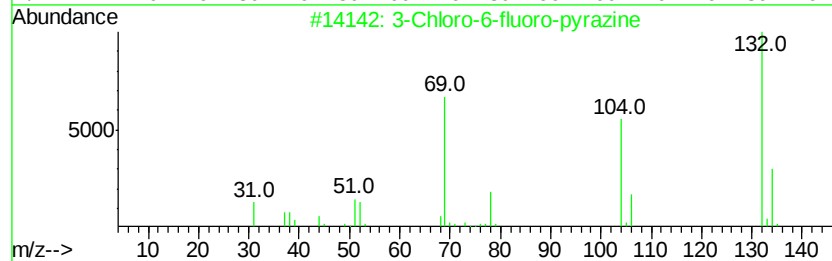
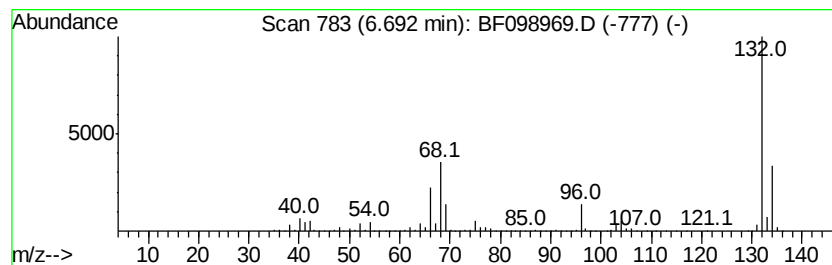
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	66.80 ng	3546960	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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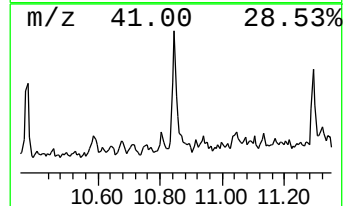
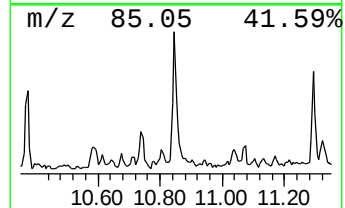
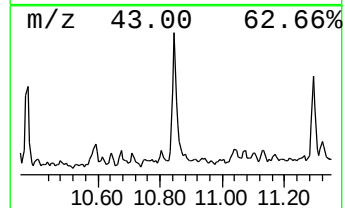
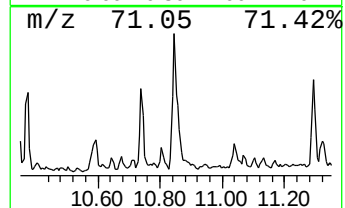
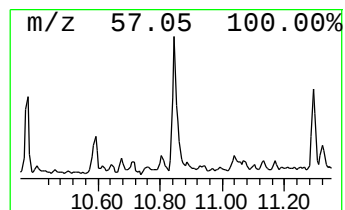
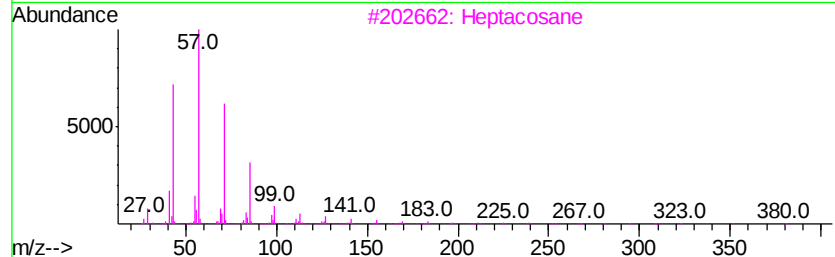
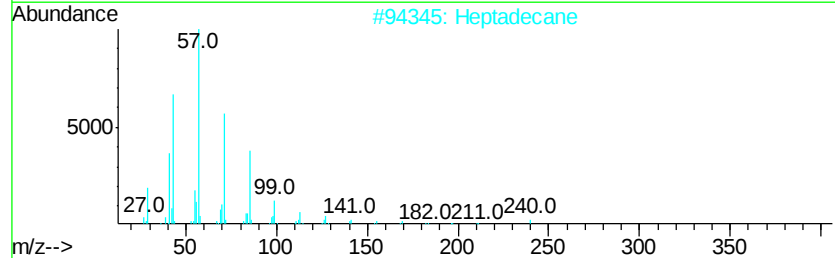
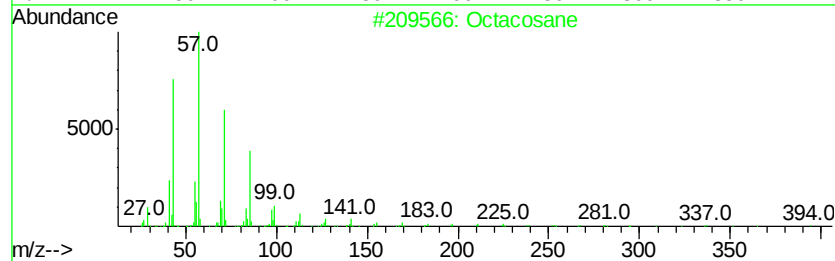
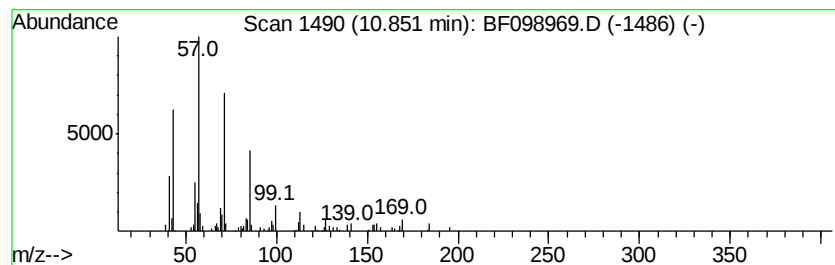
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	3.15 ng	174349	Phenanthrene-d10	11.45

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	90
2	Heptadecane		240	C17H36	000629-78-7	90
3	Heptacosane		380	C27H56	000593-49-7	90
4	Heneicosane		296	C21H44	000629-94-7	90
5	Pentadecane		212	C15H32	000629-62-9	90



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Misc :
ALS Vial : 13 Sample Multiplier: 1

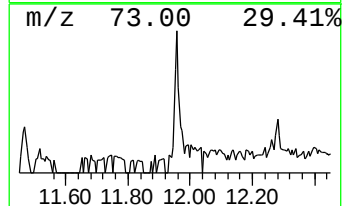
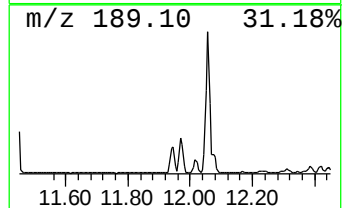
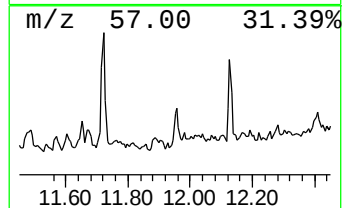
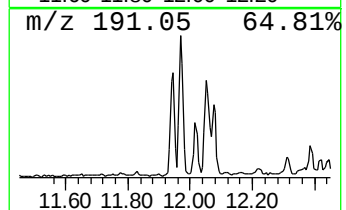
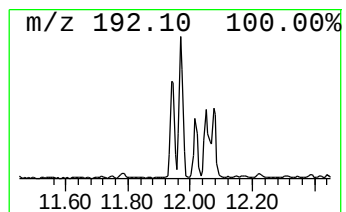
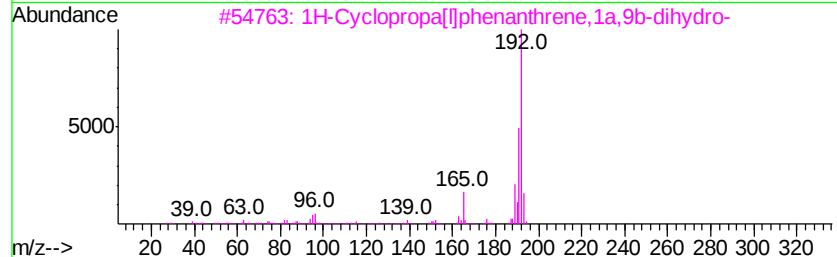
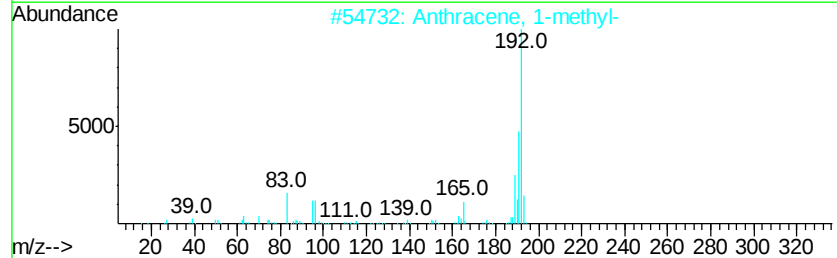
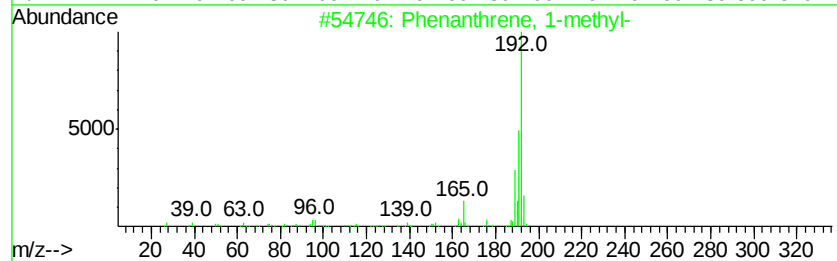
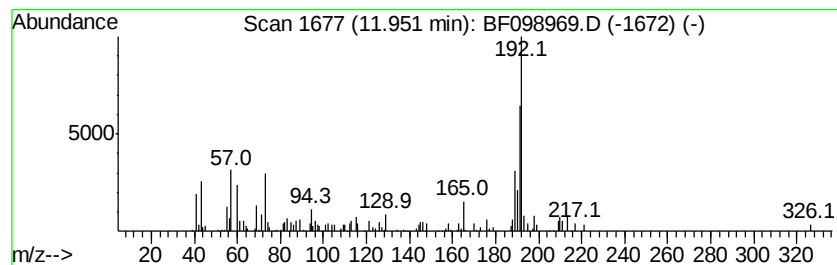
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BNA_F
ClientSampleId :
NYSEG-PH4-SB-ESMI-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.95	2.01 ng	111475	Phenanthrene-d10	11.45	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	87
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
Data File : BF098969.D
Acq On : 30 Sep 2017 18:15
Operator : SJ/JU
Sample : I5563-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

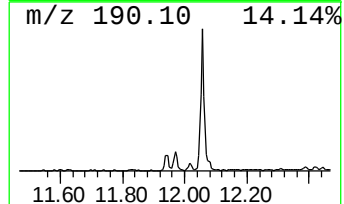
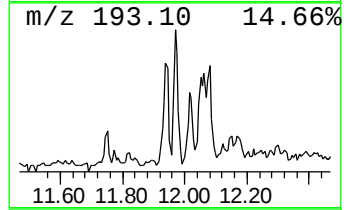
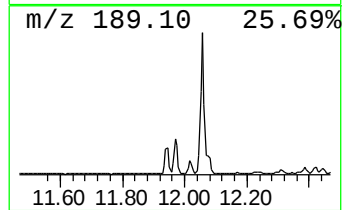
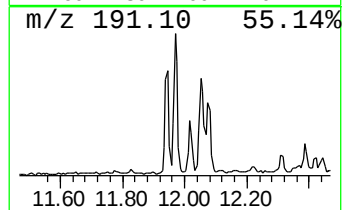
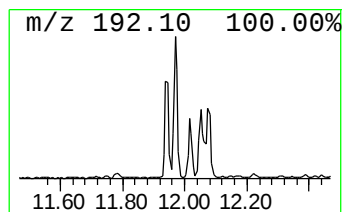
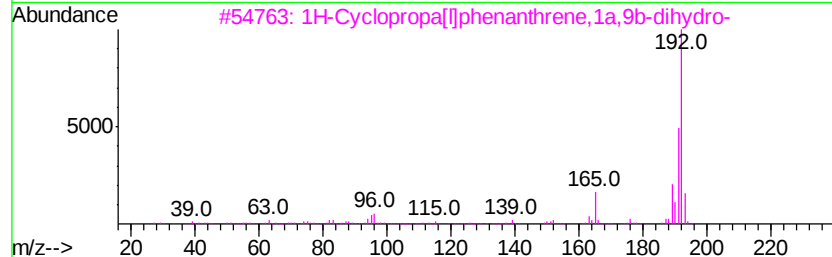
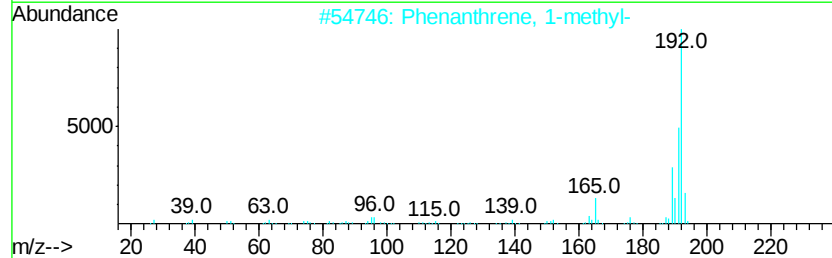
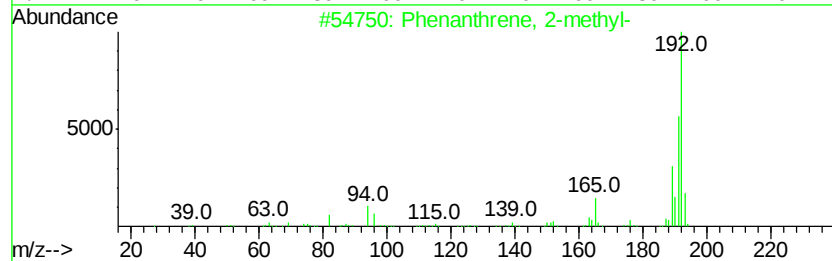
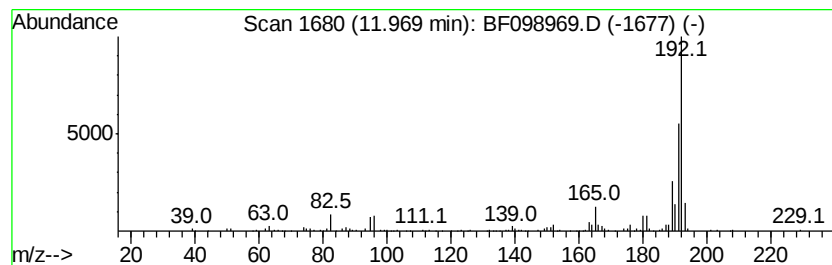
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NYSEG-PH4-SB-ESMI-6

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.97	3.30 ng	182737	Phenanthrene-d10	11.45	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
Data File : BF098969.D
Acq On : 30 Sep 2017 18:15
Operator : SJ/JU
Sample : I5563-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NYSEG-PH4-SB-ESMI-6

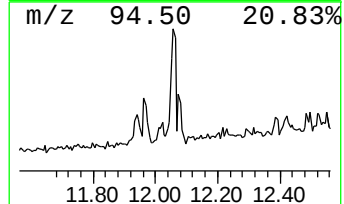
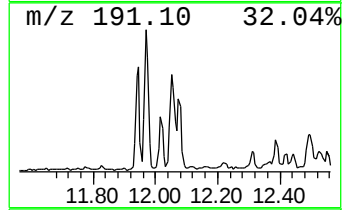
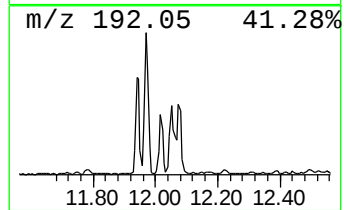
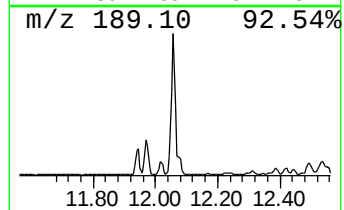
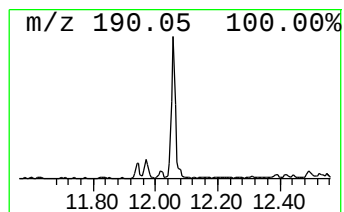
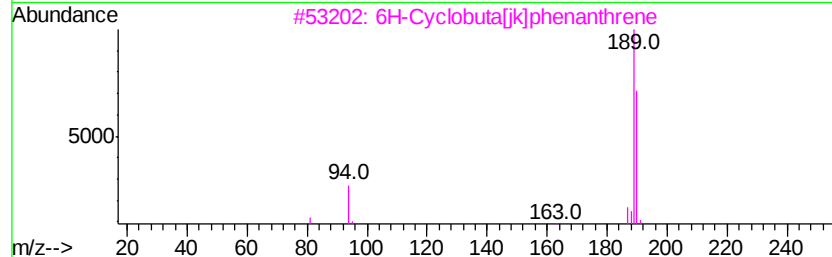
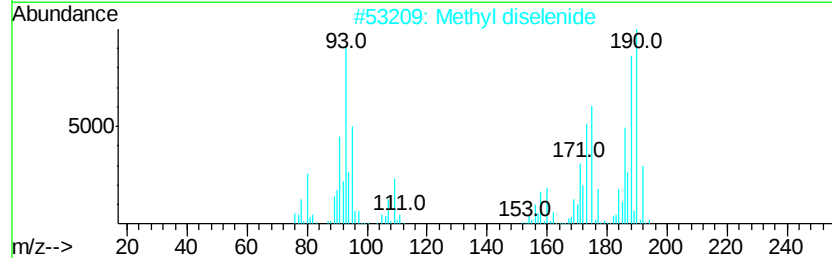
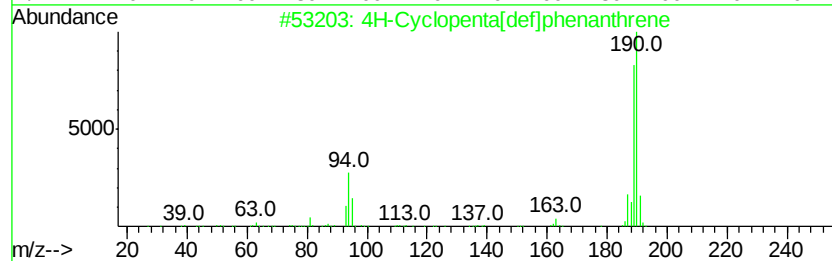
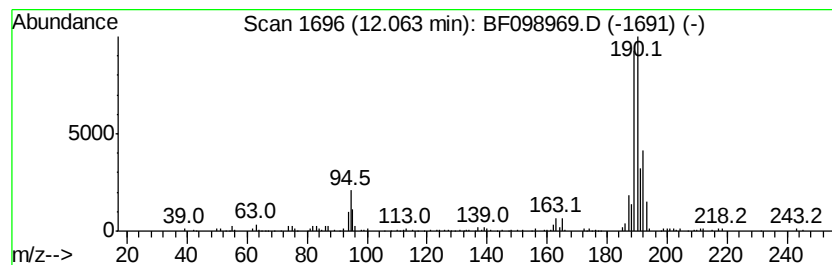
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.06	4.32 ng	239327	Phenanthrene-d10	11.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		Methyl diselenide	190	C2H6Se2	007101-31-7	55
3		6H-Cyclobuta[f]phenanthrene	190	C15H10	083469-43-6	50
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5		2-Naphthaldehyde, 1,2,3,4-tetra...	190	C12H14O2	002472-02-8	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
Data File : BF098969.D
Acq On : 30 Sep 2017 18:15
Operator : SJ/JU
Sample : I5563-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NYSEG-PH4-SB-ESMI-6

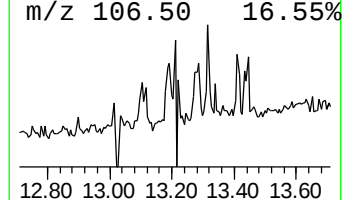
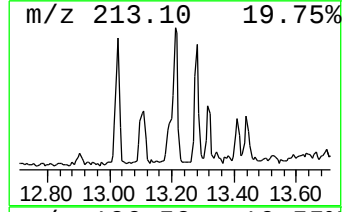
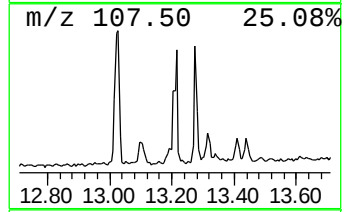
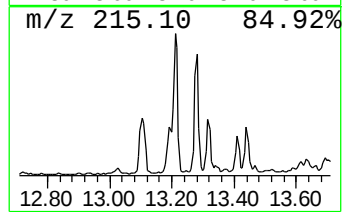
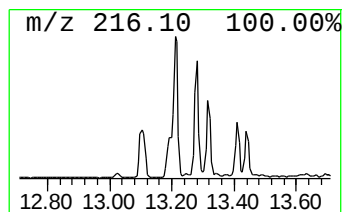
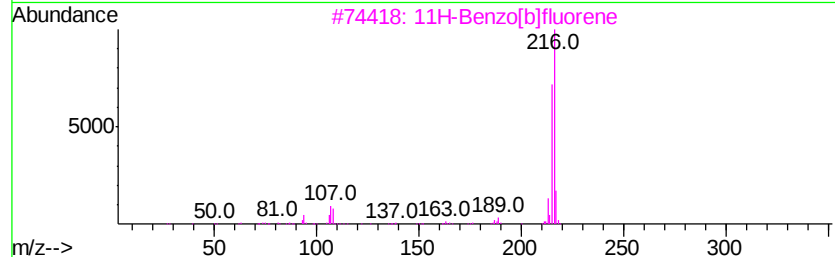
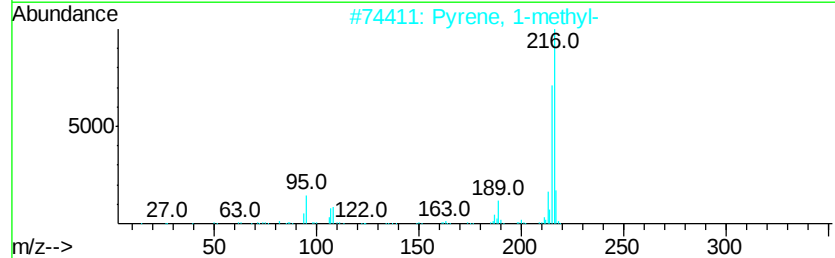
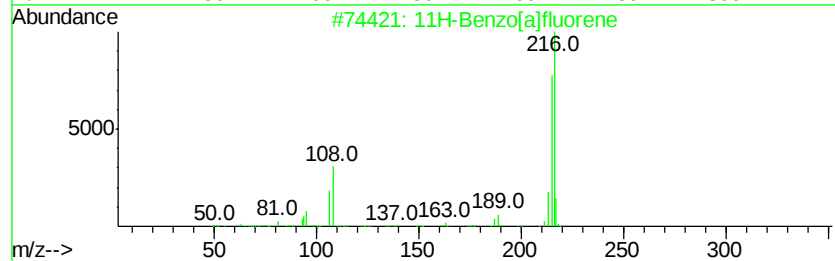
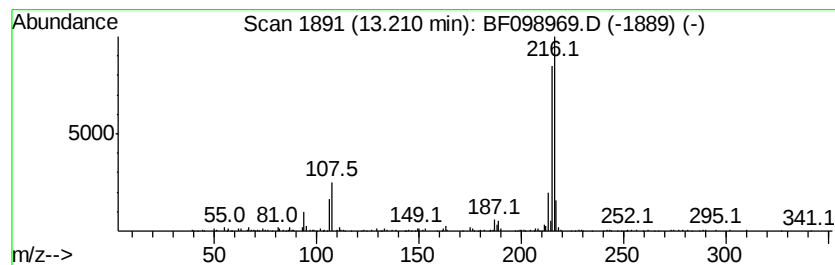
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11H-Benzo[a]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	2.27 ng	183821	Chrysene-d12	14.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
Data File : BF098969.D
Acq On : 30 Sep 2017 18:15
Operator : SJ/JU
Sample : I5563-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

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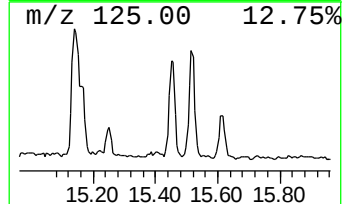
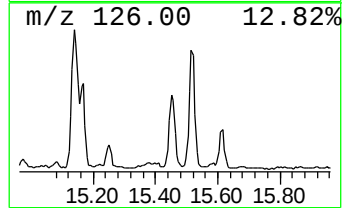
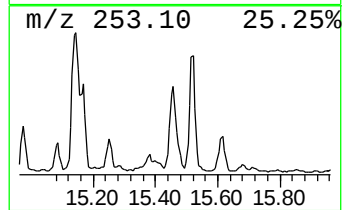
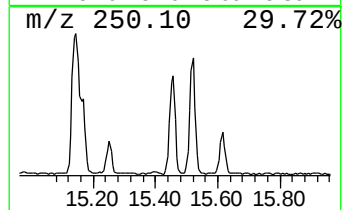
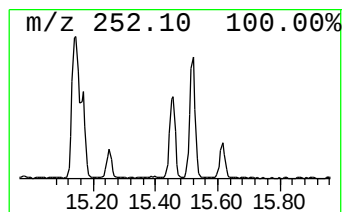
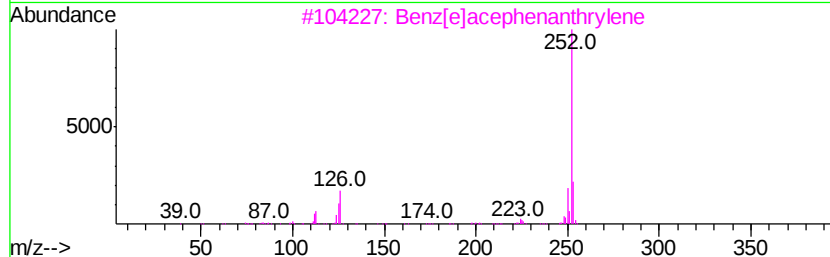
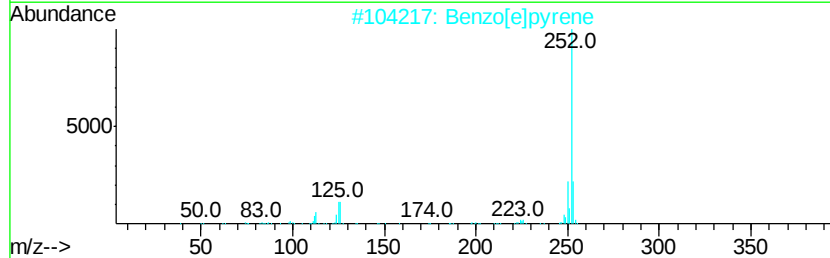
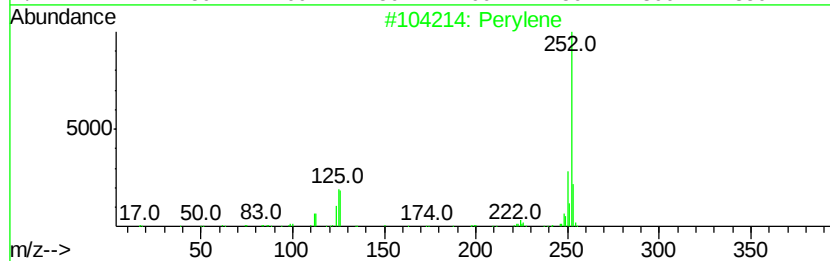
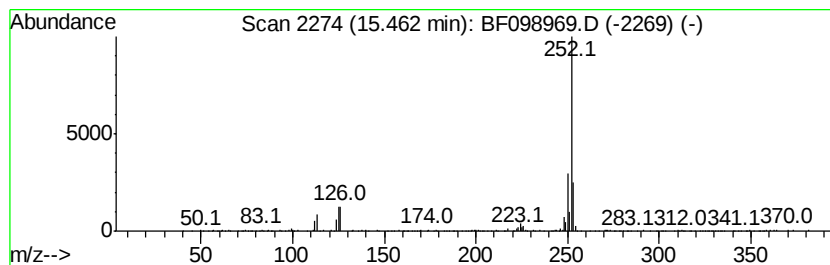
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Perylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	11.96 ng	444797	Perylene-d12	15.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Perylene	252	C20H12	000198-55-0	98
2		Benzo[e]pyrene	252	C20H12	000192-97-2	98
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	97
4		Benzo[a]pyrene	252	C20H12	000050-32-8	96
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
Data File : BF098969.D
Acq On : 30 Sep 2017 18:15
Operator : SJ/JU
Sample : I5563-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NYSEG-PH4-SB-ESMI-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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...	2.23	4.6	ng	245872	1	6.92	1061940	20.0
2-Pentanone, 4-hy...	5.15	11.1	ng	590765	1	6.92	1061940	20.0
unknown6.69	6.69	66.8	ng	3546960	1	6.92	1061940	20.0
Octacosane	10.85	3.1	ng	174349	4	11.45	1108660	20.0
Phenanthrene, 1-m...	11.95	2.0	ng	111475	4	11.45	1108660	20.0
Phenanthrene, 2-m...	11.97	3.3	ng	182737	4	11.45	1108660	20.0
4H-Cyclopenta[def...	12.06	4.3	ng	239327	4	11.45	1108660	20.0
11H-Benzo[a]fluorene	13.21	2.3	ng	183821	5	14.09	1620720	20.0
Perylene	15.46	12.0	ng	444797	6	15.59	743958	20.0