

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113514.D
 Acq On : 2 Apr 2019 19:16
 Operator : JU/SJ
 Sample : K2180-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7

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-BF032519.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	4.875	469	474	488	rBV	657423	956222	10.09%	1.914%
2	5.310	544	548	566	rBV	3138798	3190431	33.65%	6.385%
3	6.345	719	724	740	rBV	3247944	3509616	37.02%	7.024%
4	6.469	740	745	748	rBV	3398506	3356935	35.41%	6.718%
5	6.692	779	783	792	rBV	880565	762906	8.05%	1.527%
6	6.851	805	810	813	rBV	2697518	2590075	27.32%	5.184%
7	7.257	874	879	882	rBV	2331381	2105891	22.21%	4.215%
8	7.975	997	1001	1012	rBV	1007468	984794	10.39%	1.971%
9	8.745	1128	1132	1149	rBV	543595	758969	8.01%	1.519%
10	9.045	1179	1183	1186	rBV	4265466	3866803	40.78%	7.739%
11	9.439	1247	1250	1254	rBV	275554	220486	2.33%	0.441%
12	9.716	1294	1297	1301	rBV	1086925	1034672	10.91%	2.071%
13	10.510	1427	1432	1438	rBV	2420032	2310895	24.37%	4.625%
14	11.198	1545	1549	1551	rBV	1158001	983654	10.37%	1.969%
15	11.221	1551	1553	1557	rVV	940184	737385	7.78%	1.476%
16	11.274	1559	1562	1566	rVB	182028	166479	1.76%	0.333%
17	11.433	1583	1589	1591	rBV2	82049	95920	1.01%	0.192%
18	11.698	1630	1634	1637	rBV	129412	124045	1.31%	0.248%
19	11.733	1637	1640	1644	rVV2	293530	315874	3.33%	0.632%
20	11.810	1650	1653	1655	rBV	190545	192188	2.03%	0.385%
21	12.286	1730	1734	1739	rVB4	65207	107992	1.14%	0.216%
22	12.410	1751	1755	1758	rBV	1220897	1127410	11.89%	2.256%
23	12.633	1787	1793	1796	rBV	1181713	1171250	12.35%	2.344%
24	12.780	1814	1818	1821	rBV	3280494	2845887	30.02%	5.696%
25	12.857	1825	1831	1834	rBV3	86754	138975	1.47%	0.278%
26	12.939	1842	1845	1847	rBV3	82242	100086	1.06%	0.200%
27	12.963	1847	1849	1852	rVB	155833	136902	1.44%	0.274%
28	13.027	1856	1860	1863	rVV3	107333	110210	1.16%	0.221%
29	13.068	1863	1867	1874	rVB3	119222	180734	1.91%	0.362%
30	13.157	1879	1882	1885	rBV	98832	94930	1.00%	0.190%
31	13.357	1910	1916	1919	rBV4	58086	103817	1.09%	0.208%
32	13.474	1933	1936	1940	rVB	118373	123092	1.30%	0.246%
33	13.580	1950	1954	1956	rBV	125962	136247	1.44%	0.273%
34	13.662	1966	1968	1970	rBV	156872	135858	1.43%	0.272%

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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-BF032519.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.827	1990	1996	1999	rBV	6957281	9481102	100.00%	18.975%
36	13.857	1999	2001	2004	rVB	575950	457889	4.83%	0.916%
37	13.998	2020	2025	2028	rVV4	87807	129653	1.37%	0.259%
38	14.221	2058	2063	2066	rBV	125858	142067	1.50%	0.284%
39	14.304	2074	2077	2081	rVB2	106969	119667	1.26%	0.239%
40	14.527	2112	2115	2123	rBV5	52725	98384	1.04%	0.197%
41	14.598	2124	2127	2130	rBV2	116361	112844	1.19%	0.226%
42	14.839	2164	2168	2176	rBV	669395	1191025	12.56%	2.384%
43	14.909	2177	2180	2183	rVV	138397	142335	1.50%	0.285%
44	14.939	2183	2185	2188	rVB	115382	109168	1.15%	0.218%
45	15.121	2212	2216	2221	rVB	351582	455215	4.80%	0.911%
46	15.180	2221	2226	2230	rBV	544334	626134	6.60%	1.253%
47	15.233	2231	2235	2238	rBV	750781	951947	10.04%	1.905%
48	15.651	2302	2306	2311	rBV	84515	109775	1.16%	0.220%
49	15.886	2341	2346	2356	rVB	167034	305346	3.22%	0.611%
50	16.586	2459	2465	2471	rBV	238154	437955	4.62%	0.876%
51	16.992	2528	2534	2543	rVB	155552	318560	3.36%	0.638%

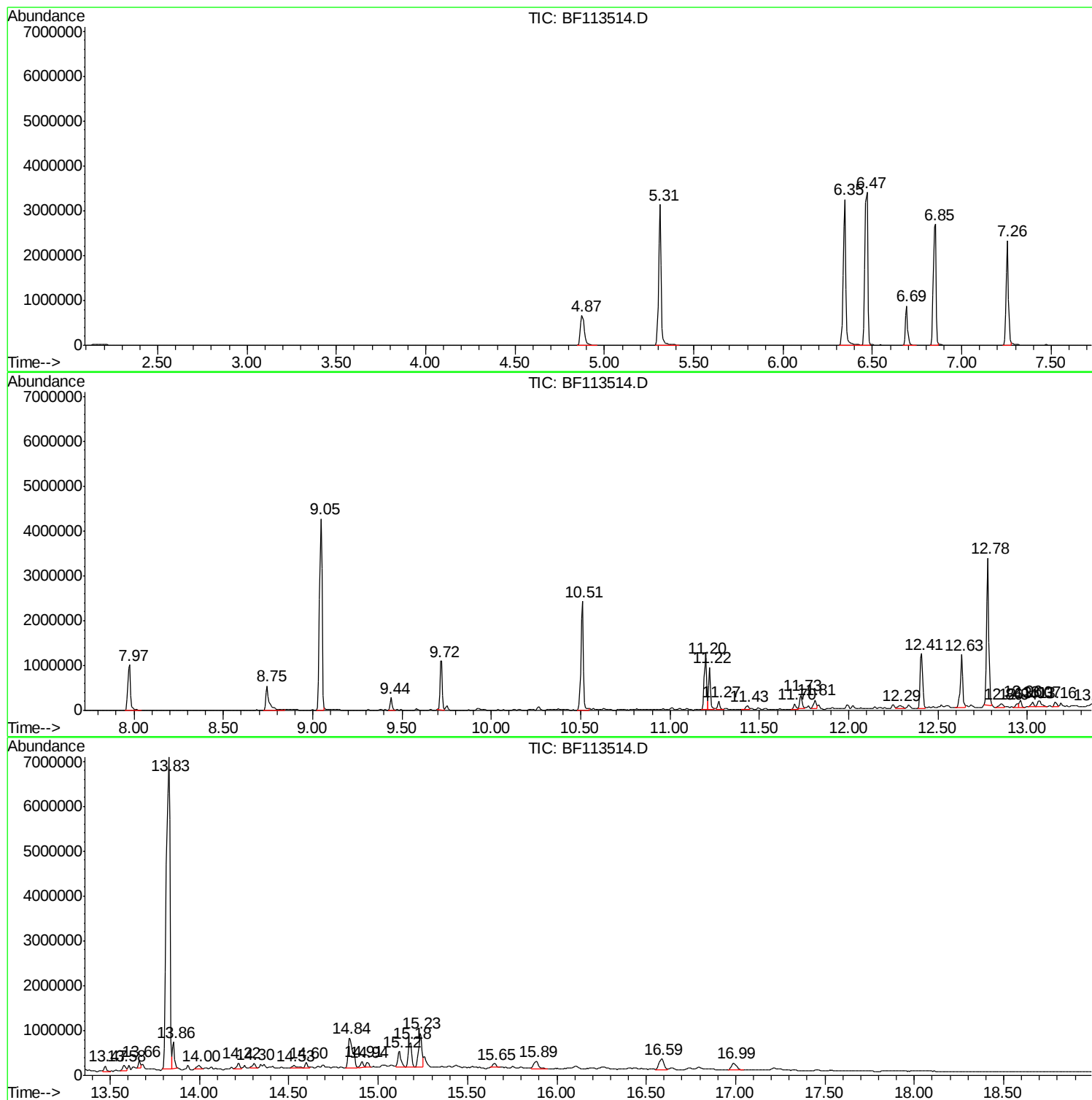
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ClientSampled :
TP-7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.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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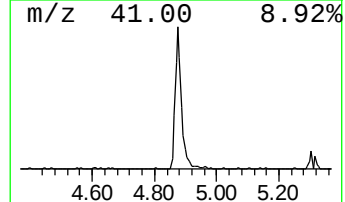
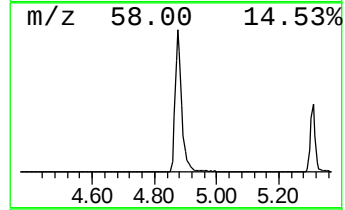
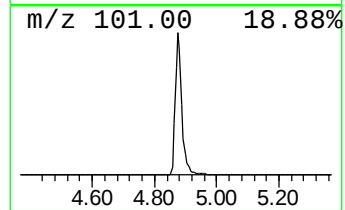
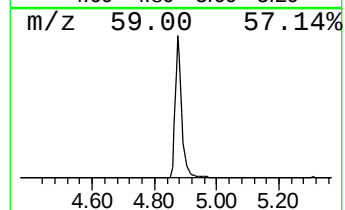
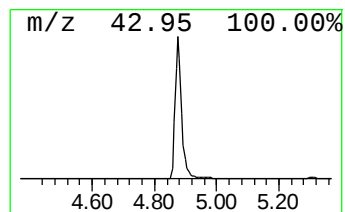
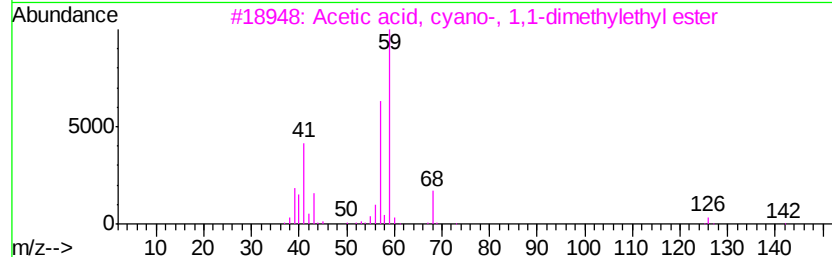
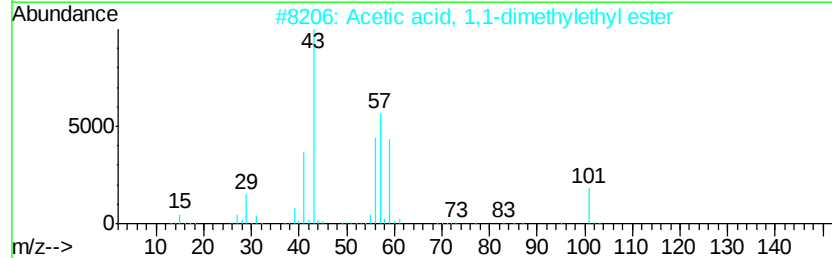
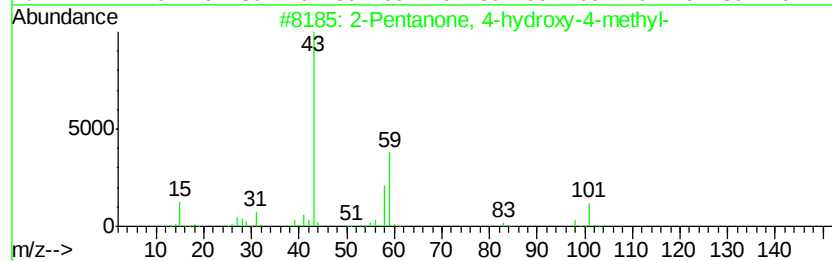
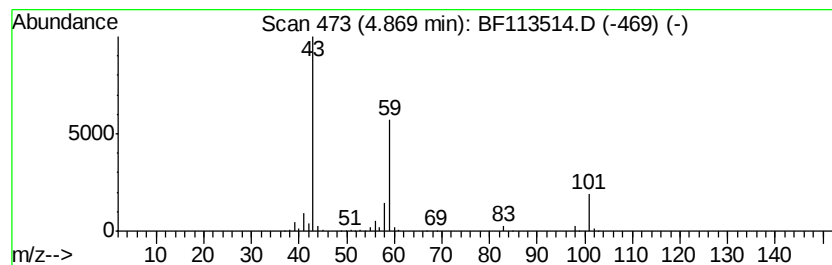
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.87	25.07 ng	956222	1,4-Dichlorobenzene-d4	6.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
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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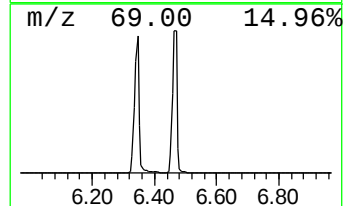
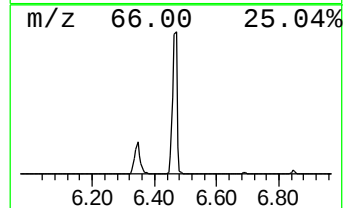
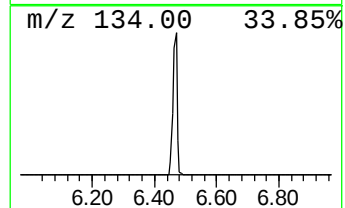
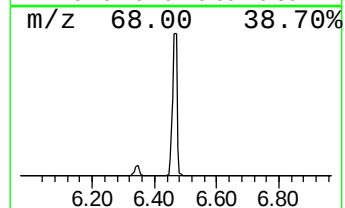
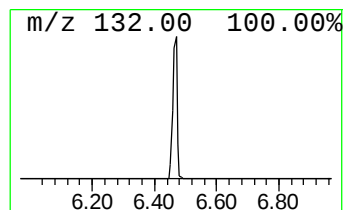
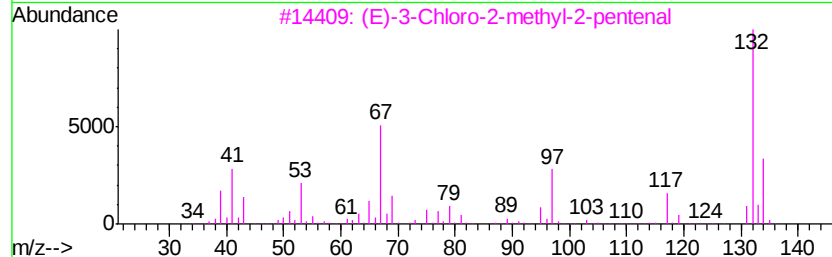
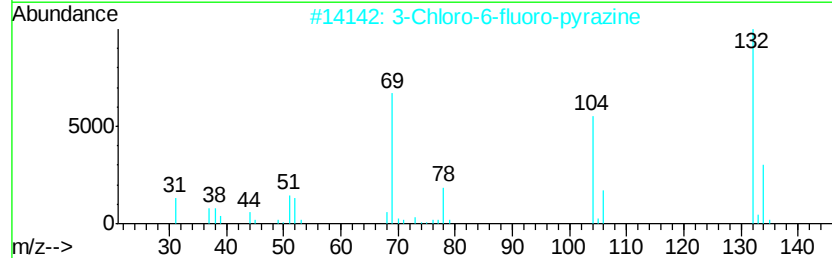
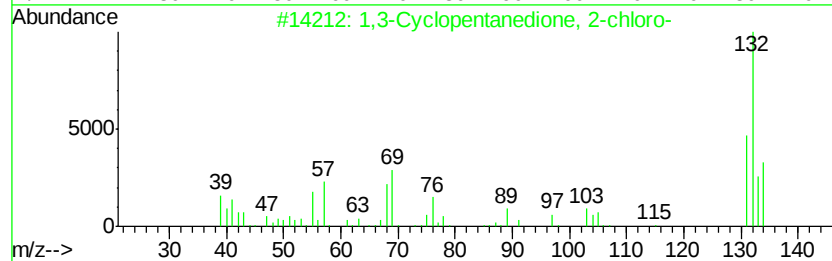
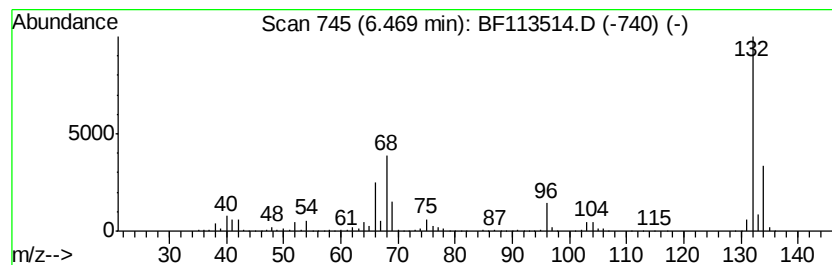
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	88.00 ng	3356940	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	47
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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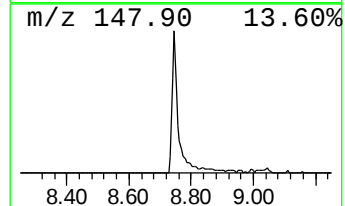
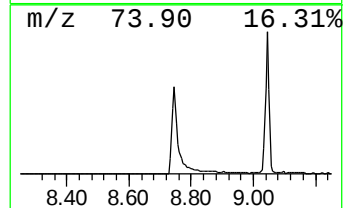
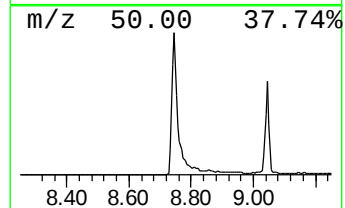
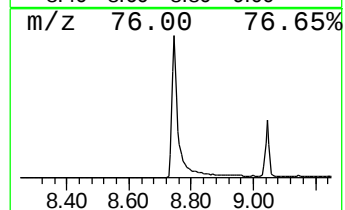
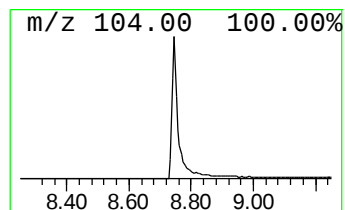
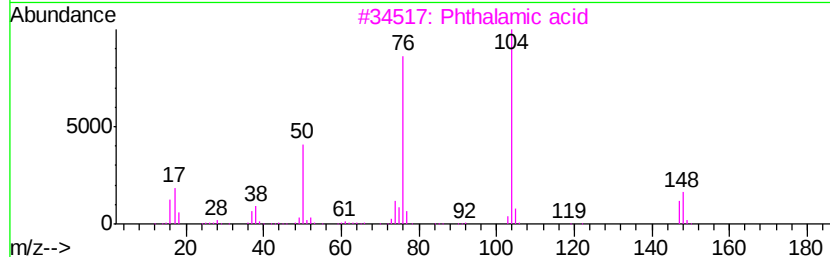
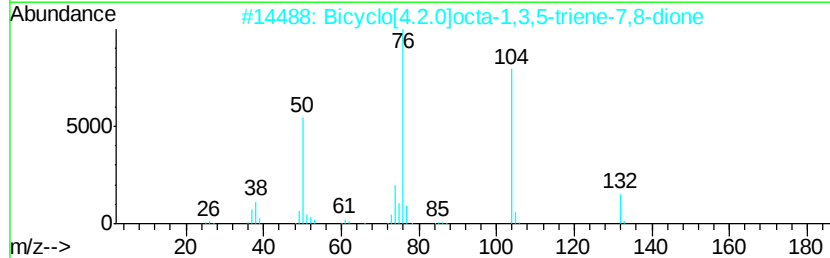
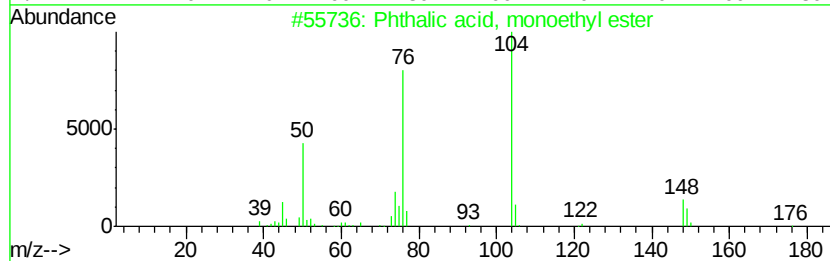
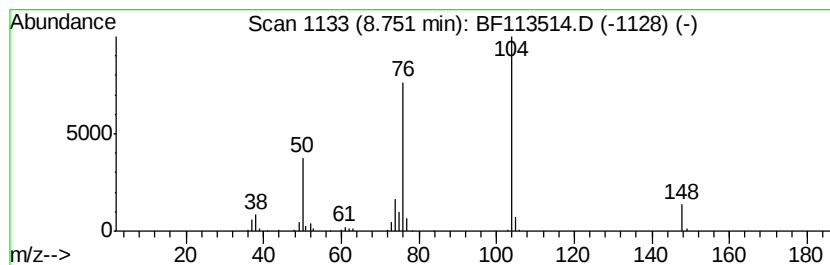
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phthalic acid, monoethyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.75	15.41 ng	758969	Naphthalene-d8	7.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, monoethyl ester	194	C10H10O4	002306-33-4	90
2		Bicyclo[4.2.0]octa-1,3,5-triene-...	132	C8H4O2	006383-11-5	86
3		Phthalamic acid	165	C8H7NO3	000088-97-1	83
4		1,2-Benzenedicarboxylic acid	166	C8H6O4	000088-99-3	83
5		Phthalic anhydride	148	C8H4O3	000085-44-9	70



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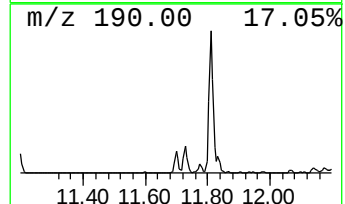
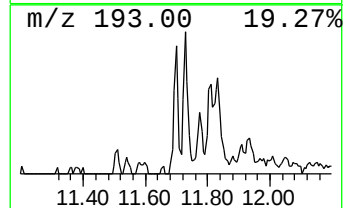
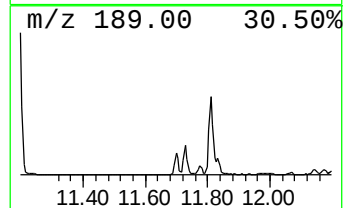
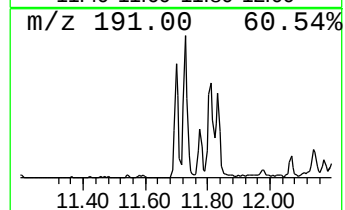
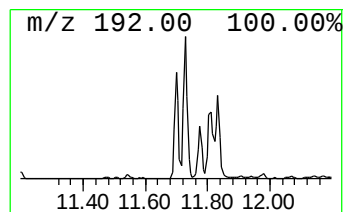
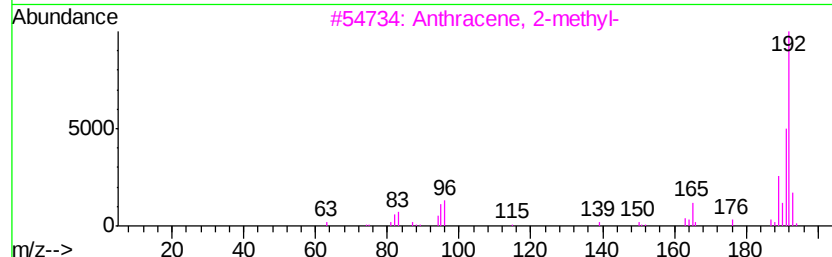
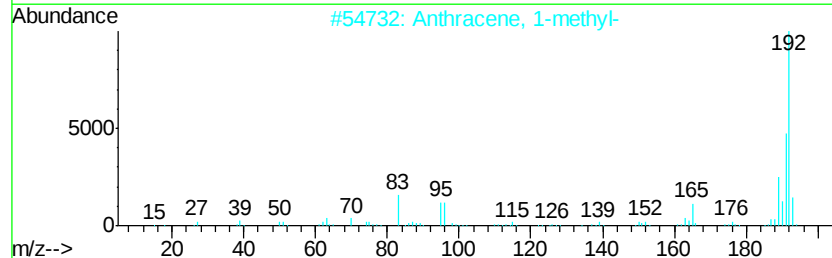
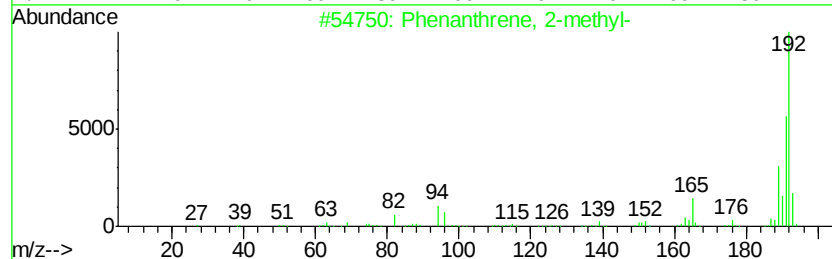
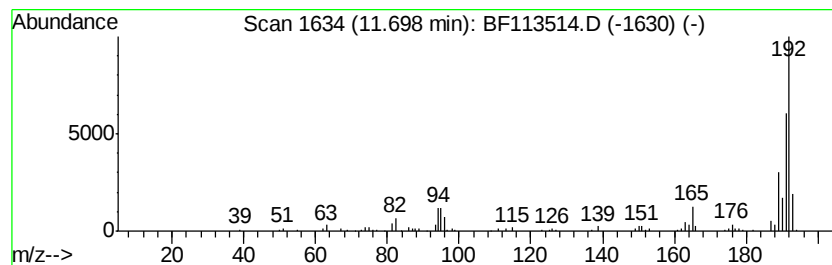
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	2.52 ng	124045	Phenanthrene-d10	11.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90



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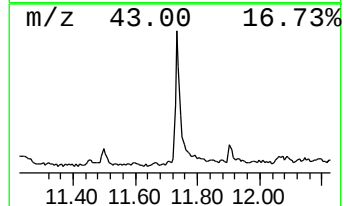
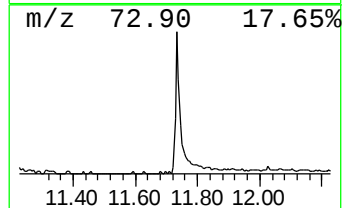
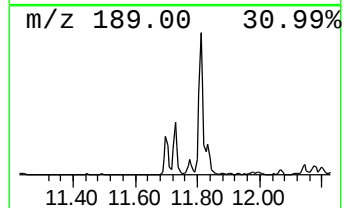
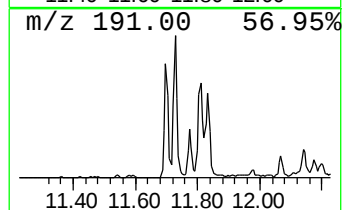
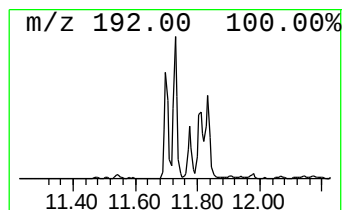
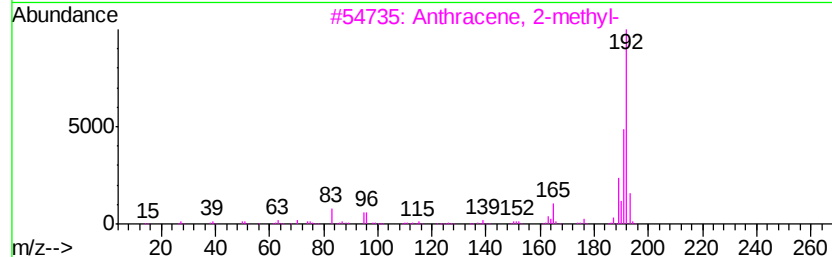
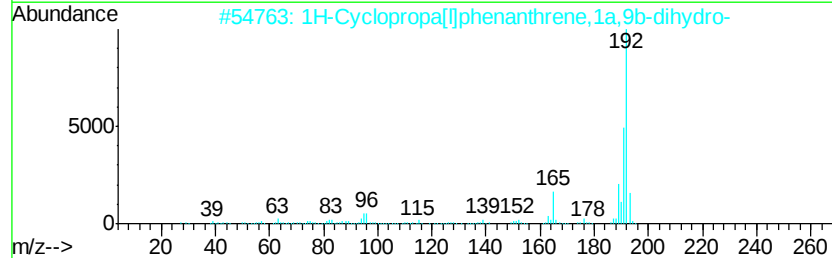
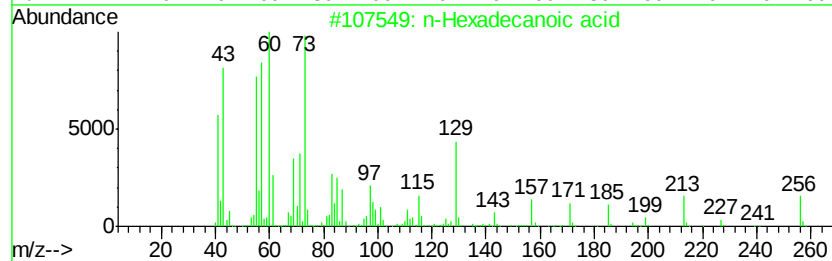
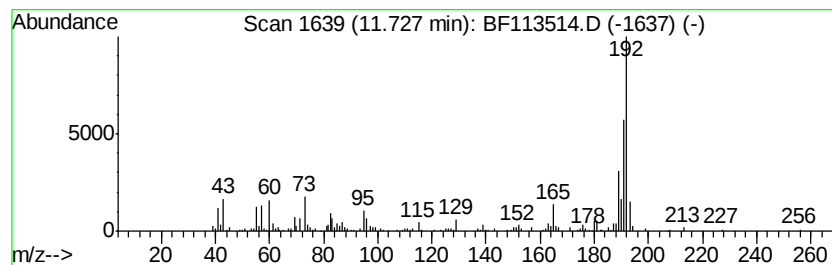
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.73	6.42 ng	315874	Phenanthrene-d10	11.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	89
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	89
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	89
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
Data File : BF113514.D
Acq On : 2 Apr 2019 19:16
Operator : JU/SJ
Sample : K2180-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-7

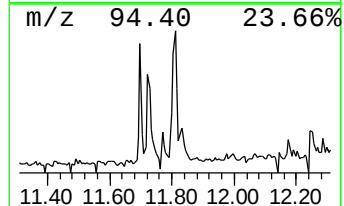
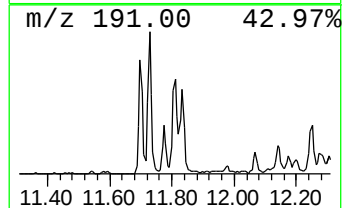
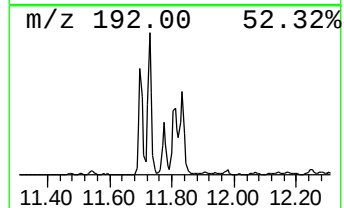
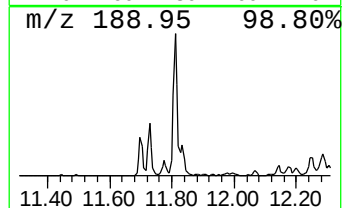
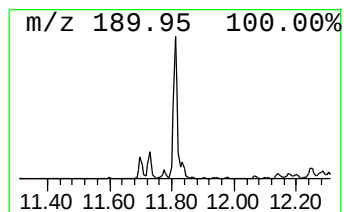
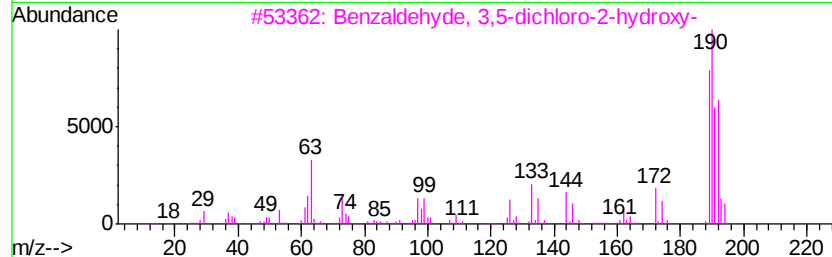
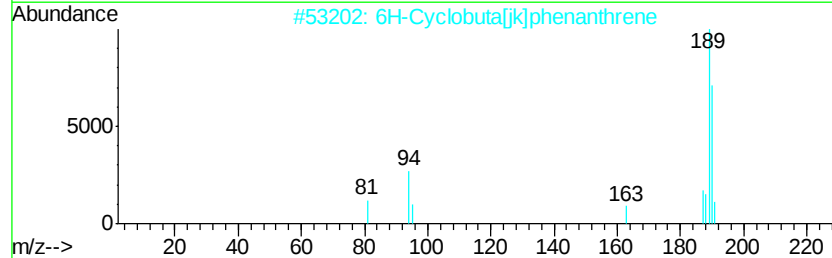
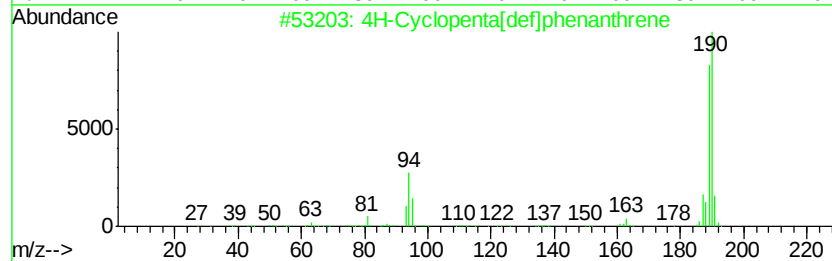
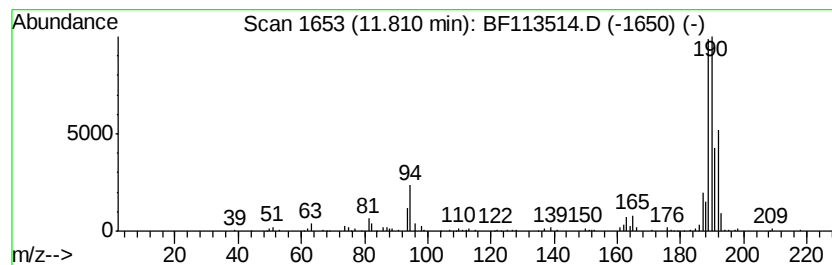
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	3.91 ng	192188	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	92
2		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
3		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
4		Methyl diselenide	190	C2H6Se2	007101-31-7	43
5		2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
Data File : BF113514.D
Acq On : 2 Apr 2019 19:16
Operator : JU/SJ
Sample : K2180-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

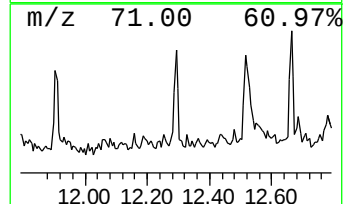
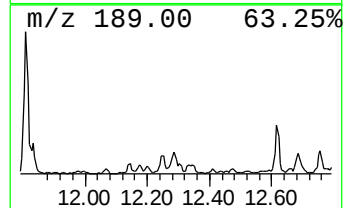
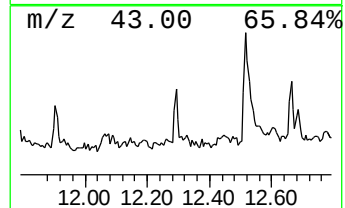
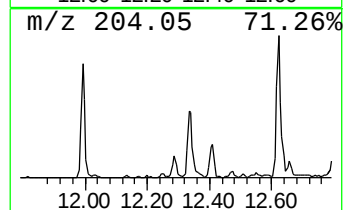
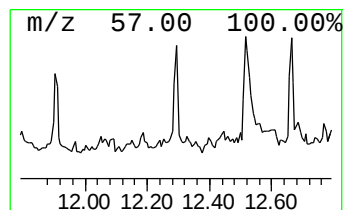
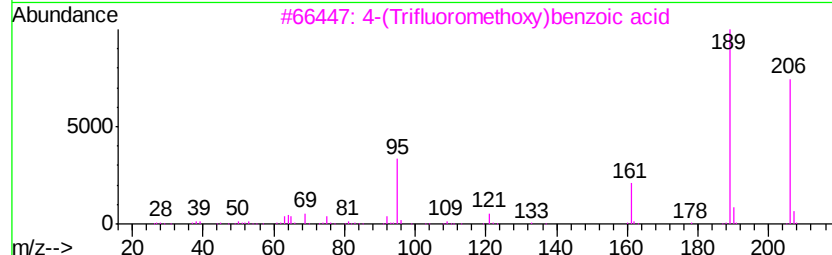
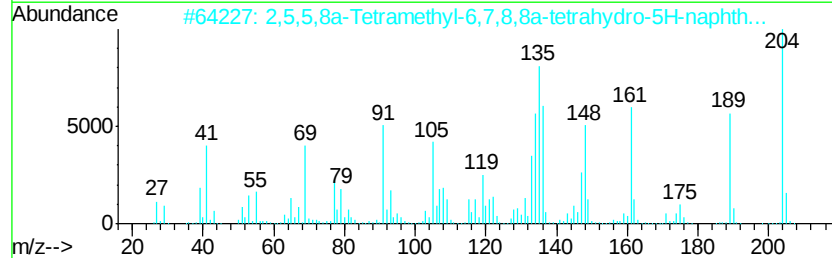
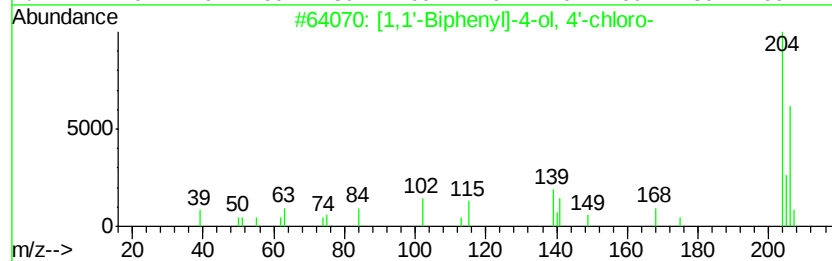
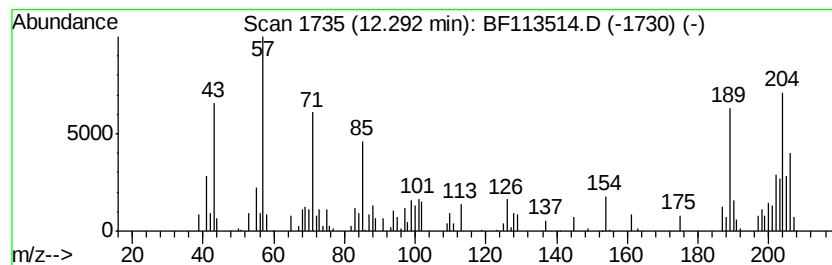
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown12.29 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.29	2.20 ng	107992	Phenanthrene-d10	11.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-ol, 4'-chloro-	204	C12H9ClO	028034-99-3	41
2	2,5,5,8a-Tetramethyl-6,7,8,8a-te...	204	C14H20O	124957-09-1	35
3	4-(Trifluoromethoxy)benzoic acid	206	C8H5F3O3	000330-12-1	22
4	Pyrene, 4,5,9,10-tetrahydro-	206	C16H14	000781-17-9	14
5	Benzoic acid, p-[[dimethylamino...	206	C11H14N2O2	029390-16-7	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
Data File : BF113514.D
Acq On : 2 Apr 2019 19:16
Operator : JU/SJ
Sample : K2180-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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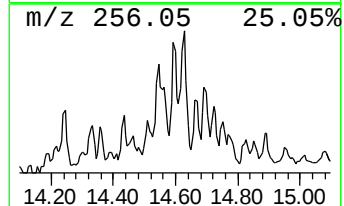
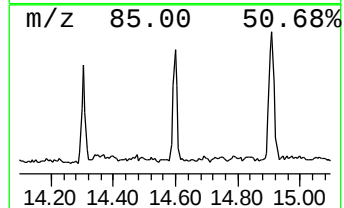
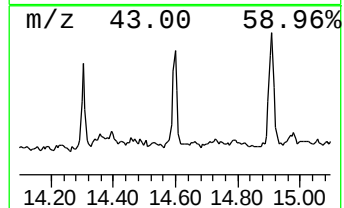
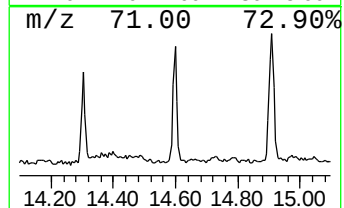
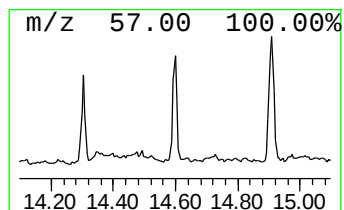
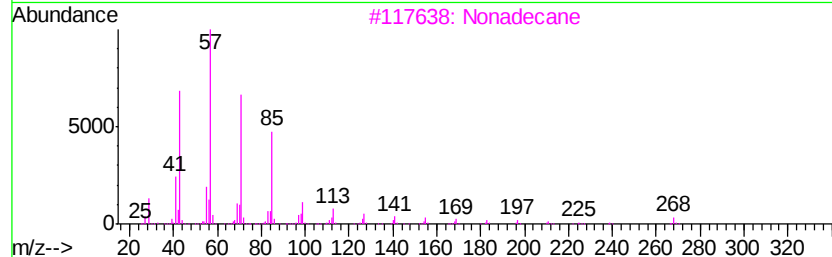
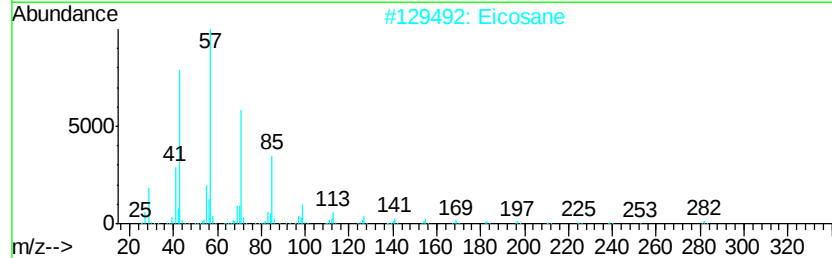
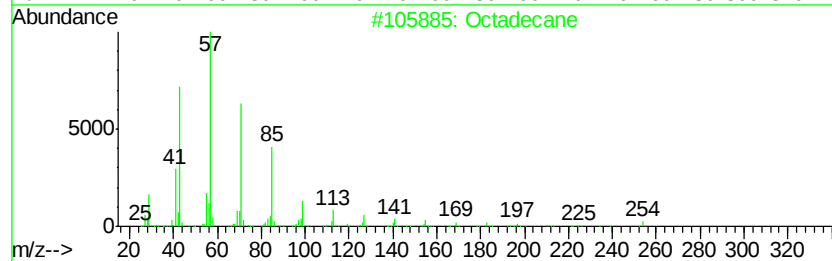
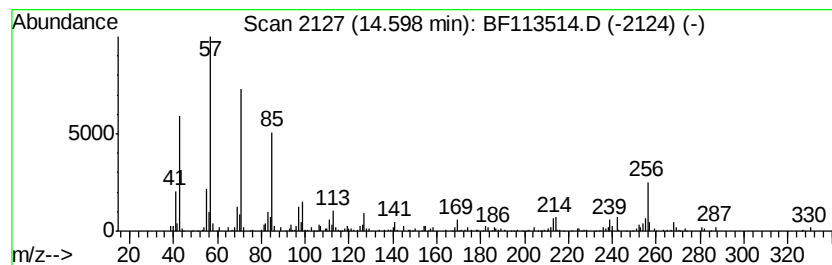
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	2.37 ng	112844	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	93
2		Eicosane	282	C20H42	000112-95-8	89
3		Nonadecane	268	C19H40	000629-92-5	89
4		Heneicosane	296	C21H44	000629-94-7	76
5		Heptadecane	240	C17H36	000629-78-7	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
Data File : BF113514.D
Acq On : 2 Apr 2019 19:16
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Sample : K2180-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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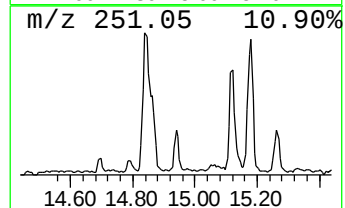
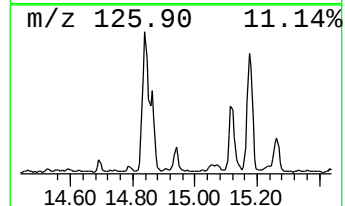
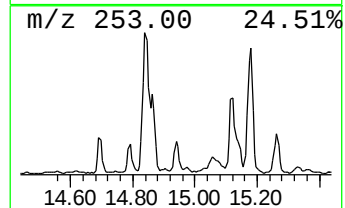
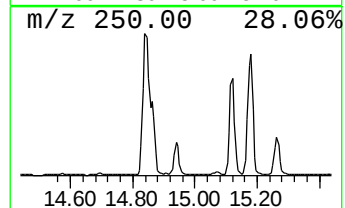
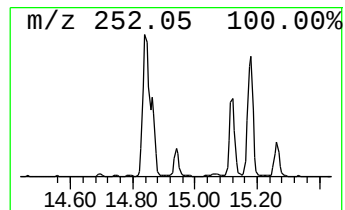
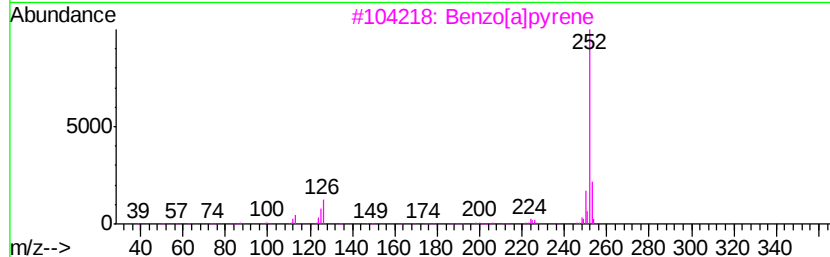
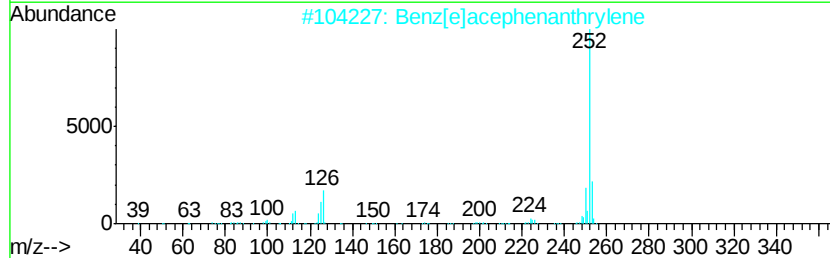
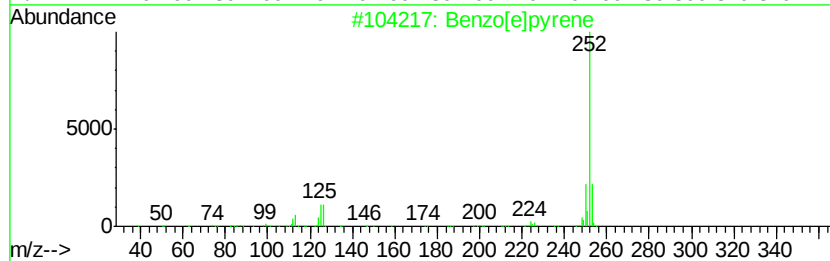
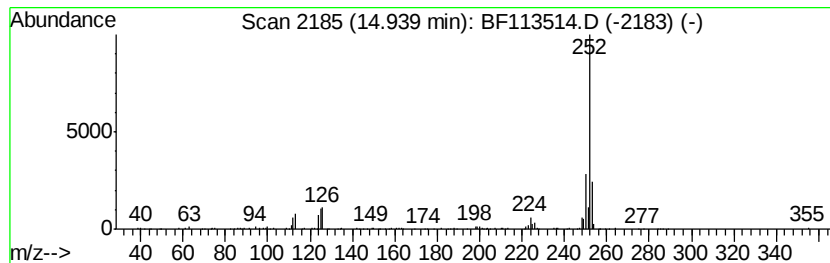
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	2.29 ng	109168	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	93
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
3		Benzo[a]pyrene	252	C20H12	000050-32-8	90
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
5		Perylene	252	C20H12	000198-55-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
Data File : BF113514.D
Acq On : 2 Apr 2019 19:16
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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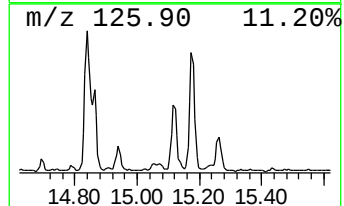
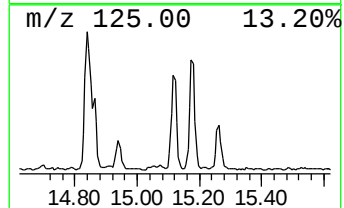
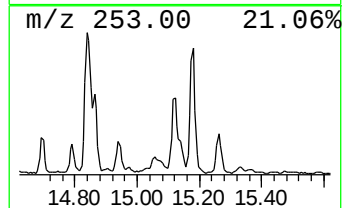
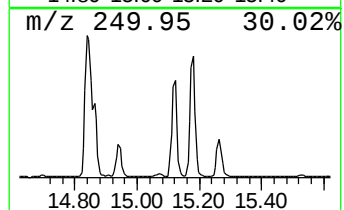
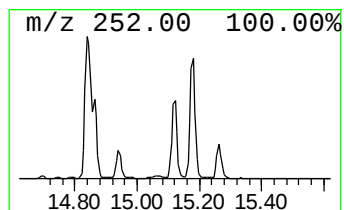
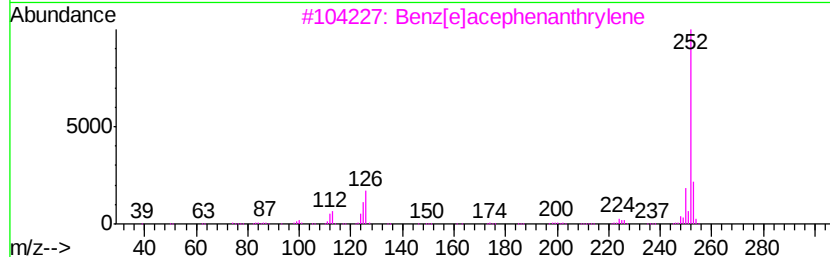
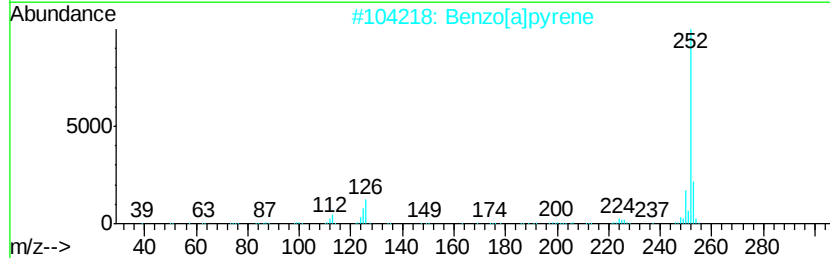
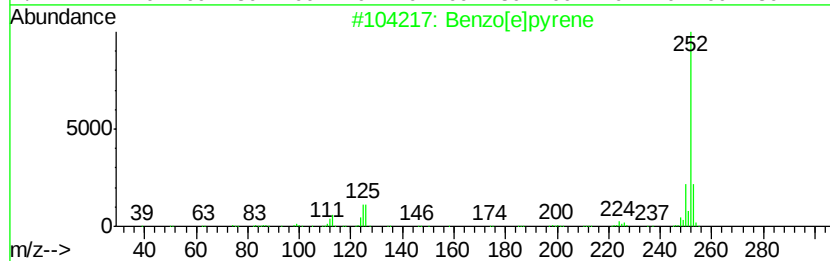
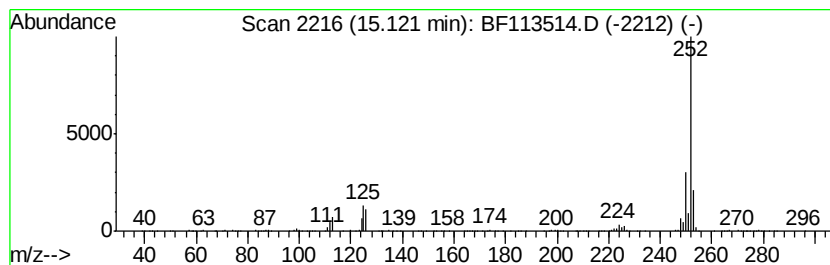
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benz[e]acephenanthrylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	9.56 ng	455215	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benzo[a]pyrene	252	C20H12	000050-32-8	98
3		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	94
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
Data File : BF113514.D
Acq On : 2 Apr 2019 19:16
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Sample : K2180-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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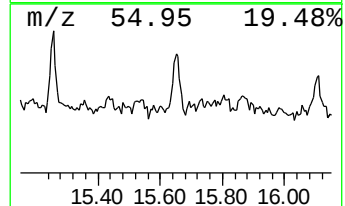
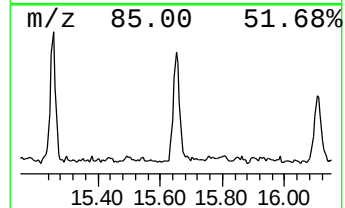
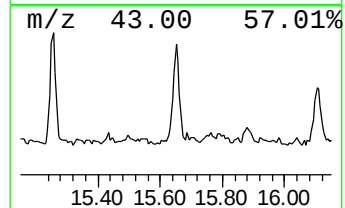
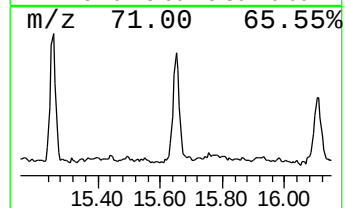
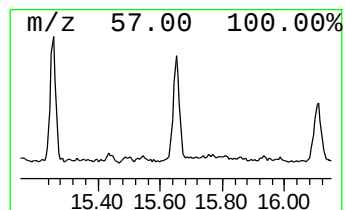
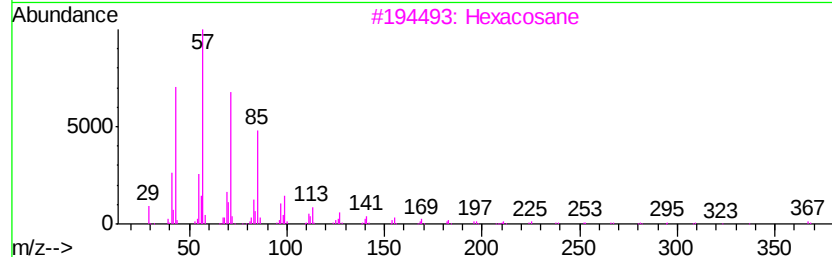
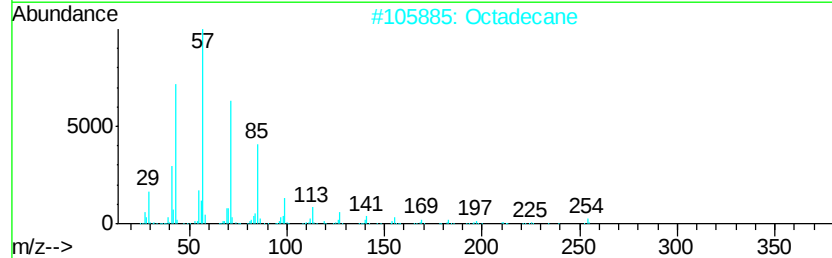
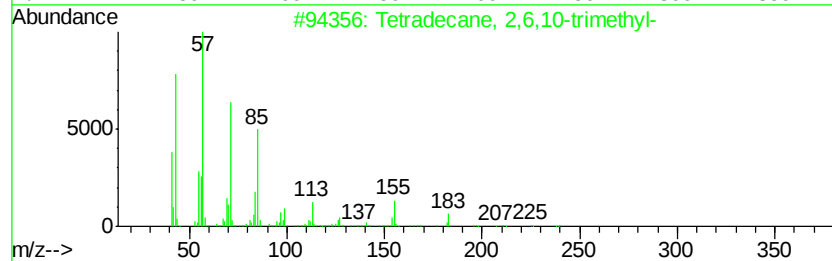
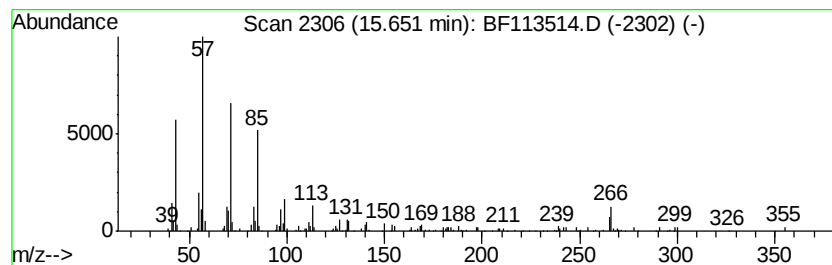
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetradecane, 2,6,10-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	2.31 ng	109775	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
2			Octadecane	254	C18H38	000593-45-3	90
3			Hexacosane	366	C26H54	000630-01-3	87
4			Pentacosane	352	C25H52	000629-99-2	83
5			Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
Data File : BF113514.D
Acq On : 2 Apr 2019 19:16
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Sample : K2180-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-7

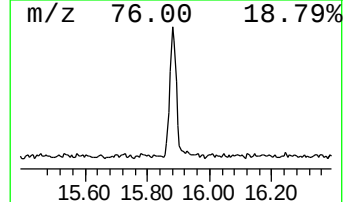
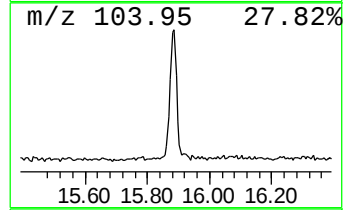
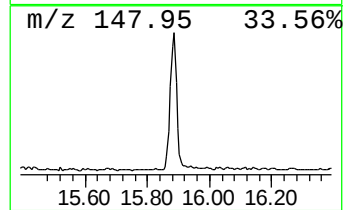
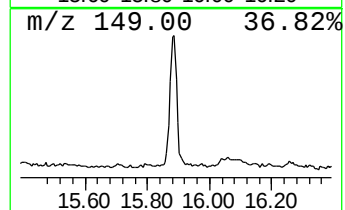
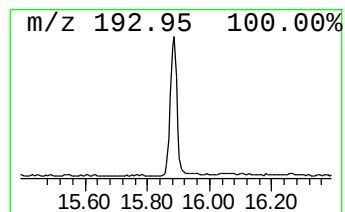
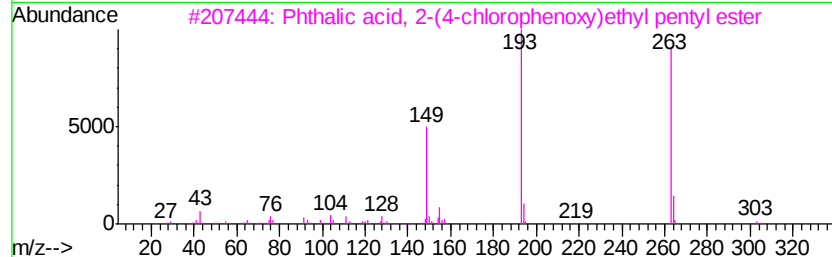
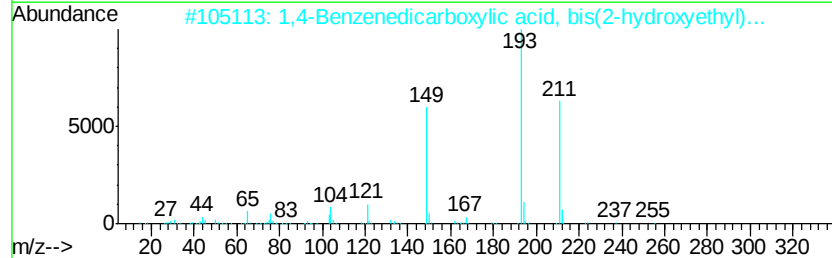
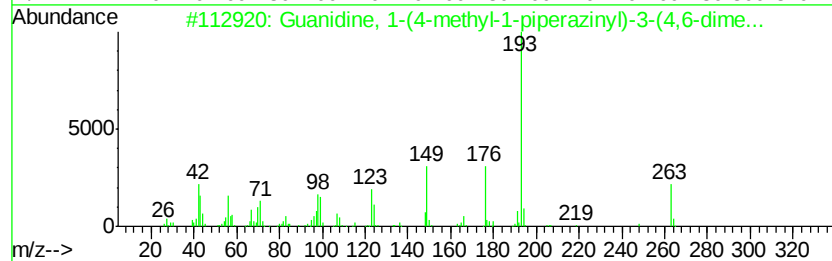
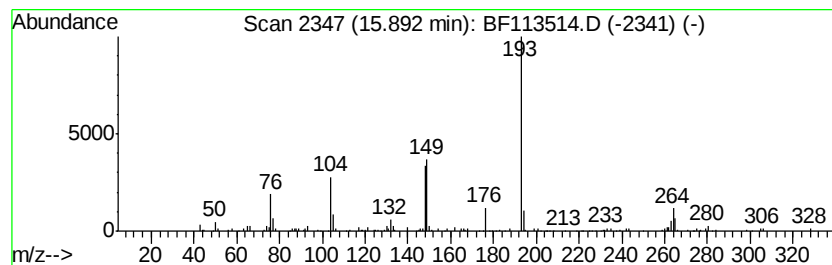
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.89 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.89	6.42 ng	305346	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Guanidine, 1-(4-methyl-1-piperaz...	263	C12H21N7	333307-74-7	33
2		1,4-Benzenedicarboxylic acid, bi...	254	C12H14O6	000959-26-2	17
3		Phthalic acid, 2-(4-chlorophenox...	390	C21H23ClO5	1000377-90-6	17
4		N-[4-Methoxy-3-methoxycarbonyl)b...	307	C14H13NO7	541528-00-1	17
5		Phthalic acid, 2-(4-chlorophenox...	376	C20H21ClO5	1000377-90-4	17



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113514.D
 Acq On : 2 Apr 2019 19:16
 Operator : JU/SJ
 Sample : K2180-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.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
2-Pentanone, 4-hy...	4.87	25.1 ng		956222	1	6.69	762906	20.0
unknown6.47	6.47	88.0 ng		3356940	1	6.69	762906	20.0
Phthalic acid, mo...	8.75	15.4 ng		758969	2	7.97	984794	20.0
Phenanthrene, 2-m...	11.70	2.5 ng		124045	4	11.20	983654	20.0
n-Hexadecanoic acid	11.73	6.4 ng		315874	4	11.20	983654	20.0
4H-Cyclopenta[def...	11.81	3.9 ng		192188	4	11.20	983654	20.0
unknown12.29	12.29	2.2 ng		107992	4	11.20	983654	20.0
Octadecane	14.60	2.4 ng		112844	6	15.23	951947	20.0
Benzo[e]pyrene	14.94	2.3 ng		109168	6	15.23	951947	20.0
Benz[e]acephenant...	15.12	9.6 ng		455215	6	15.23	951947	20.0
Tetradecane, 2,6,...	15.65	2.3 ng		109775	6	15.23	951947	20.0
unknown15.89	15.89	6.4 ng		305346	6	15.23	951947	20.0