

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122917\
 Data File : BF101837.D
 Acq On : 29 Dec 2017 18:53
 Operator : SJ/JU
 Sample : I7034-08
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TB-05(5-6)

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:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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	5.010	494	497	516	rBV	116397	240912	7.34%	0.810%
2	5.410	561	565	594	rBV	2361385	3282768	100.00%	11.044%
3	6.434	736	739	757	rBV	2429235	3187822	97.11%	10.724%
4	6.557	757	760	766	rVV	2635762	2951416	89.91%	9.929%
5	6.787	795	799	808	rBV	778643	825474	25.15%	2.777%
6	6.940	821	825	839	rBV	2352886	2327588	70.90%	7.830%
7	7.357	892	896	915	rBV	1785662	1960403	59.72%	6.595%
8	8.069	1013	1017	1035	rVB	861562	1045461	31.85%	3.517%
9	9.139	1195	1199	1207	rBV	2767039	2527713	77.00%	8.504%
10	9.204	1207	1210	1214	rVB	131211	107251	3.27%	0.361%
11	9.534	1260	1266	1273	rBV2	244569	363792	11.08%	1.224%
12	9.734	1297	1300	1304	rVB	211678	163524	4.98%	0.550%
13	9.822	1311	1315	1319	rBV	909025	835209	25.44%	2.810%
14	9.851	1319	1320	1328	rVB	48806	63147	1.92%	0.212%
15	9.963	1336	1339	1342	rBV4	26048	36793	1.12%	0.124%
16	10.051	1352	1354	1359	rVB2	46526	61851	1.88%	0.208%
17	10.233	1378	1385	1388	rBV	235942	214908	6.55%	0.723%
18	10.381	1407	1410	1415	rVB	33646	41856	1.28%	0.141%
19	10.451	1417	1422	1425	rBV	89279	101984	3.11%	0.343%
20	10.622	1447	1451	1462	rBV	1223582	1408689	42.91%	4.739%
21	10.704	1462	1465	1470	rBV	185806	225369	6.87%	0.758%
22	11.151	1538	1541	1544	rBV2	89999	79956	2.44%	0.269%
23	11.180	1544	1546	1548	rVB	57872	42828	1.30%	0.144%
24	11.316	1565	1569	1571	rBV	871531	785831	23.94%	2.644%
25	11.339	1571	1573	1580	rVB	356221	328411	10.00%	1.105%
26	11.398	1580	1583	1586	rBV	86705	88410	2.69%	0.297%
27	11.580	1609	1614	1616	rBV2	64220	76012	2.32%	0.256%
28	11.828	1652	1656	1658	rBV2	205518	210786	6.42%	0.709%
29	11.933	1671	1674	1676	rBV	74805	81394	2.48%	0.274%
30	11.986	1680	1683	1686	rBV	65926	49227	1.50%	0.166%
31	12.116	1702	1705	1707	rBV	32943	36354	1.11%	0.122%
32	12.369	1745	1748	1753	rBV2	85386	114226	3.48%	0.384%
33	12.533	1773	1776	1784	rBV	554182	557952	17.00%	1.877%
34	12.610	1787	1789	1795	rBV5	43836	68928	2.10%	0.232%

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35	12.763	1810	1815	1821	rBV	546687	708998	21.60%	2.385%
36	12.898	1834	1838	1841	rBV	2064565	1968663	59.97%	6.623%
37	13.774	1984	1987	1993	rBV2	439300	507072	15.45%	1.706%
38	13.969	2016	2020	2023	rBV2	839155	1002386	30.53%	3.372%
39	13.998	2023	2025	2028	rVB	224666	191474	5.83%	0.644%
40	15.021	2196	2199	2202	rBV	278911	331729	10.11%	1.116%
41	15.451	2268	2272	2276	rBV	384340	520677	15.86%	1.752%

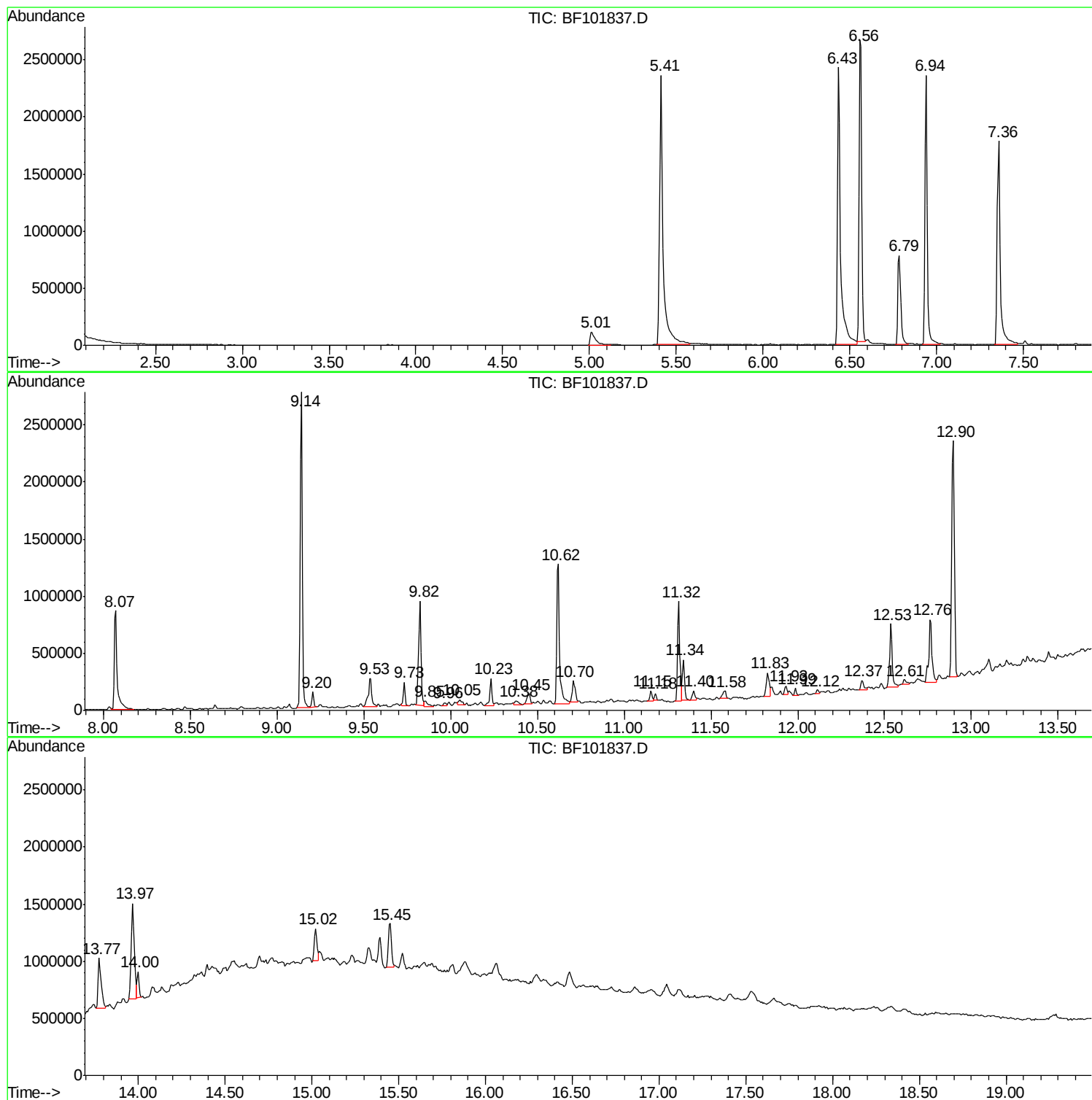
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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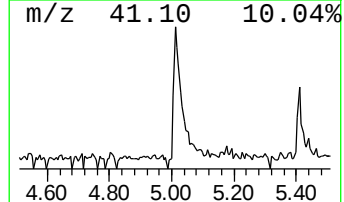
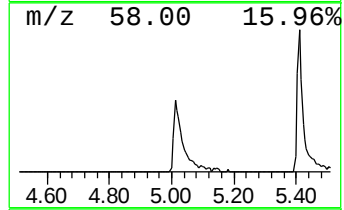
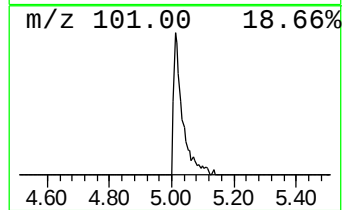
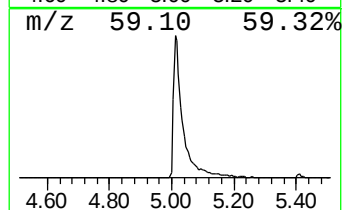
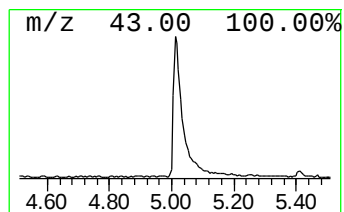
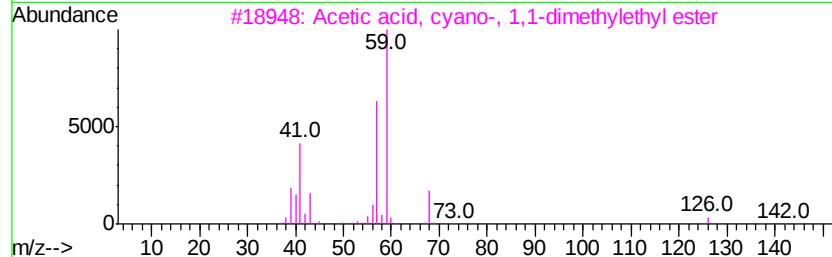
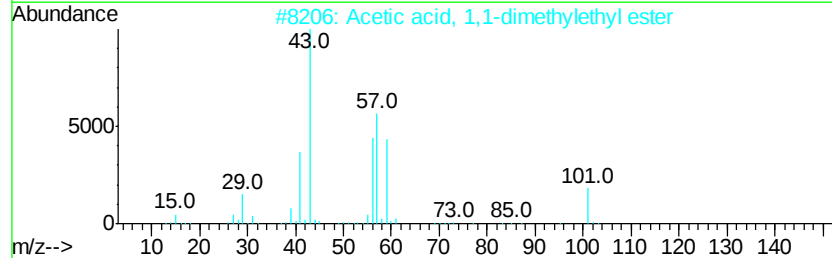
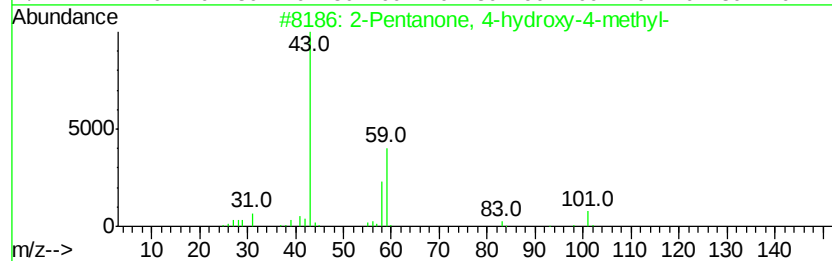
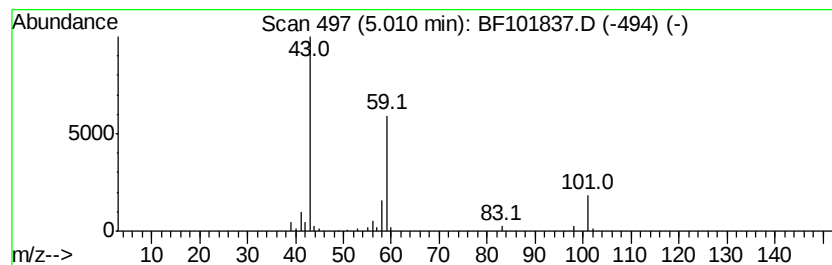
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	5.84 ng	240912	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-			116	C6H12O2	000123-42-2	45
2	Acetic acid, 1,1-dimethylethyl e...			116	C6H12O2	000540-88-5	38
3	Acetic acid, cyano-, 1,1-dimethyl...			141	C7H11NO2	001116-98-9	23
4	5-Hexen-2-one			98	C6H10O	000109-49-9	9
5	Morpholine, 4-methyl-			101	C5H11NO	000109-02-4	9



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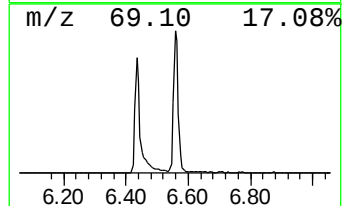
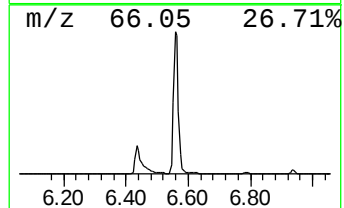
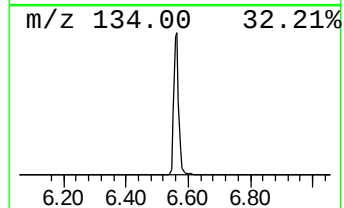
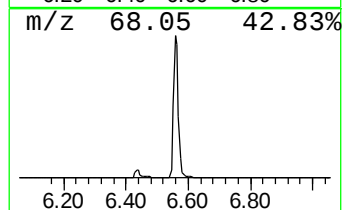
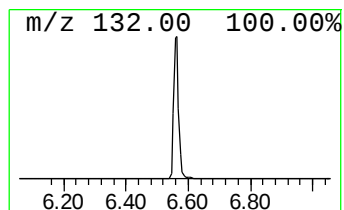
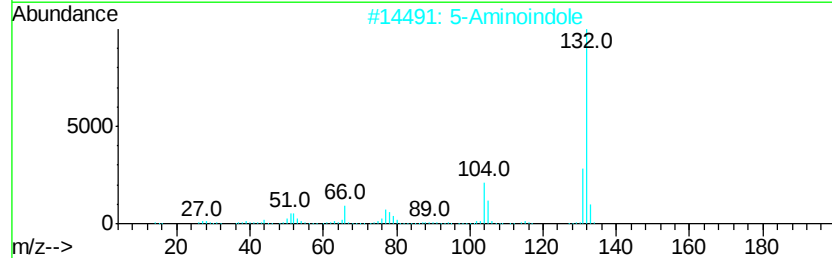
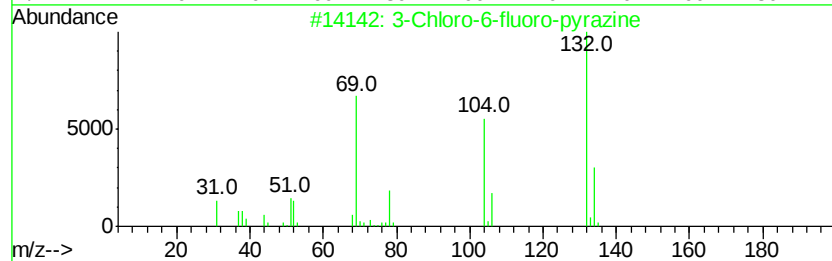
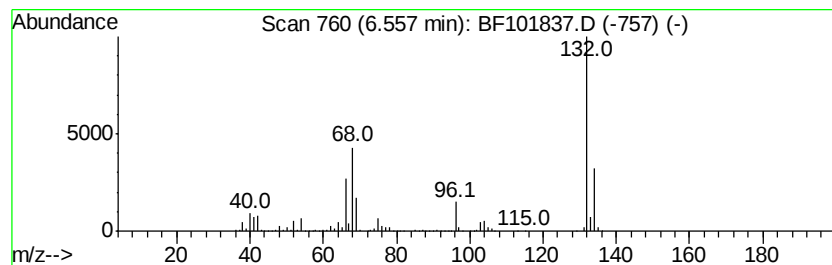
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	71.51 ng	2951420	1,4-Dichlorobenzene-d4	6.79

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



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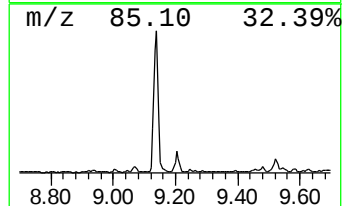
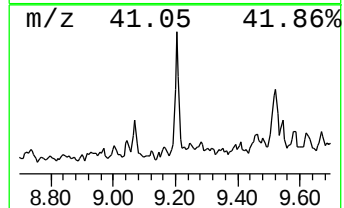
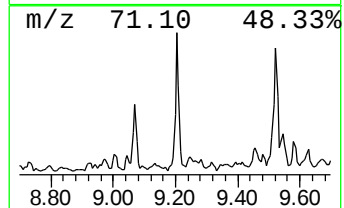
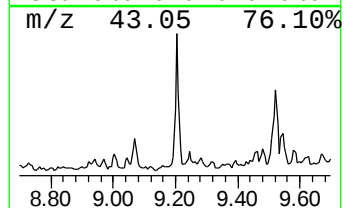
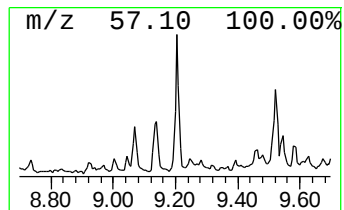
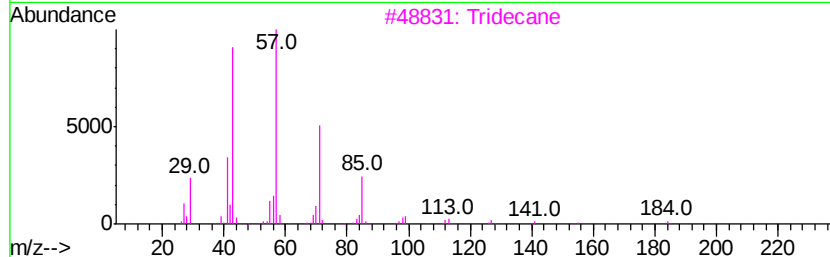
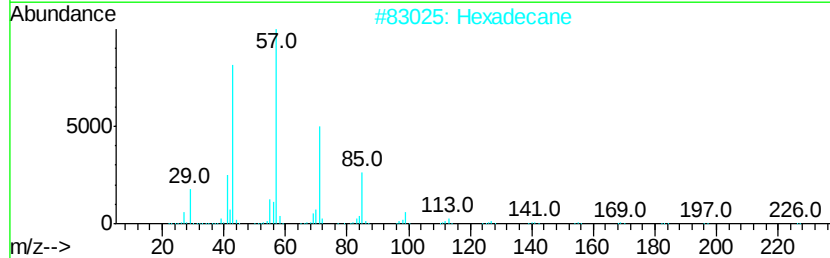
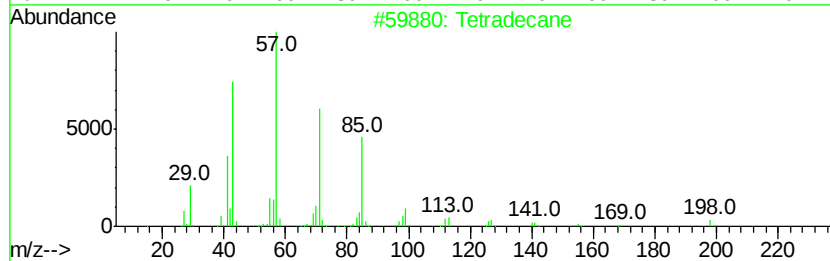
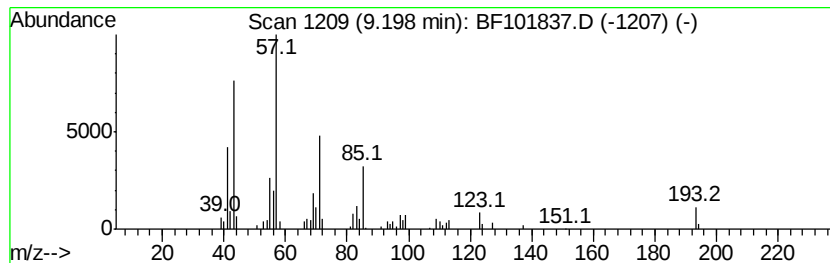
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Tetradecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.20	2.57 ng	107251	Acenaphthene-d10	9.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	80
2		Hexadecane	226	C16H34	000544-76-3	80
3		Tridecane	184	C13H28	000629-50-5	80
4		Pentadecane	212	C15H32	000629-62-9	80
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	72



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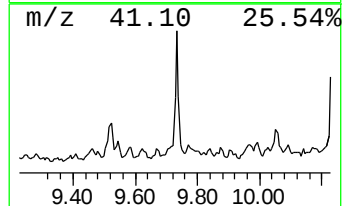
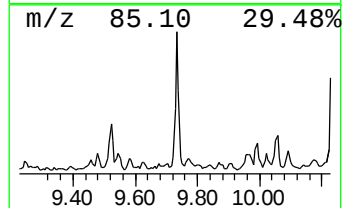
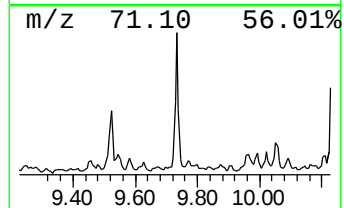
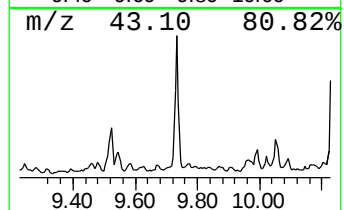
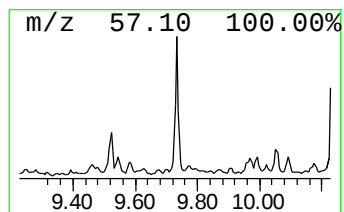
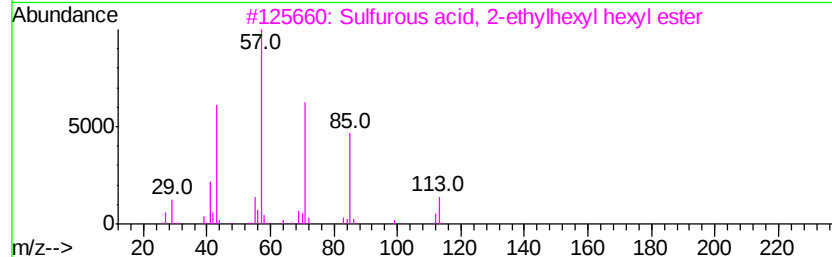
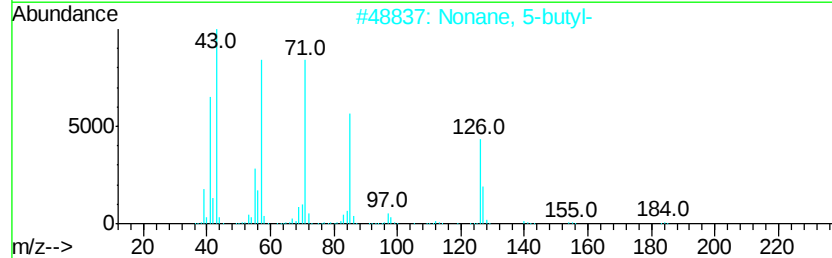
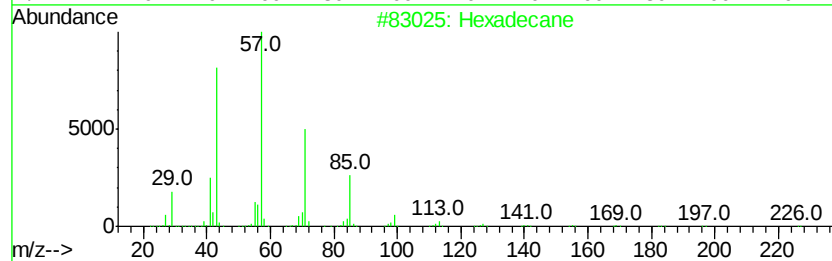
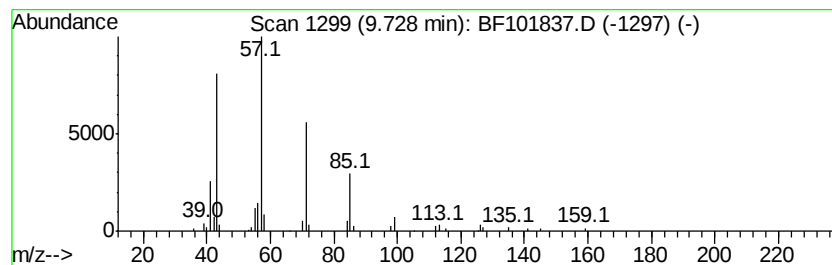
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Hexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.73	3.92 ng	163524	Acenaphthene-d10	9.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	86
2		Nonane, 5-butyl-	184	C13H28	017312-63-9	64
3		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
4		Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	59
5		Nonane, 1-iodo-	254	C9H19I	004282-42-2	50



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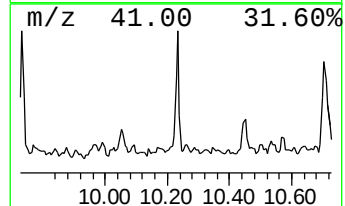
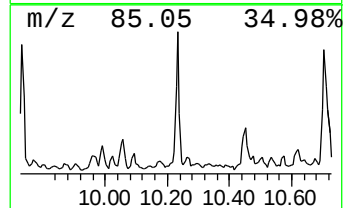
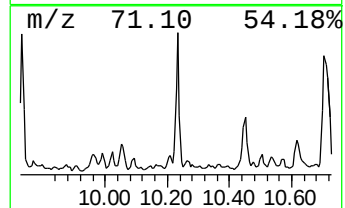
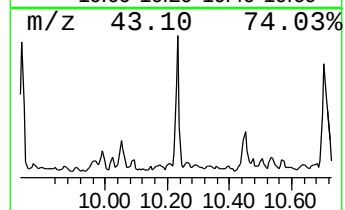
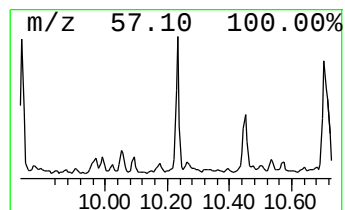
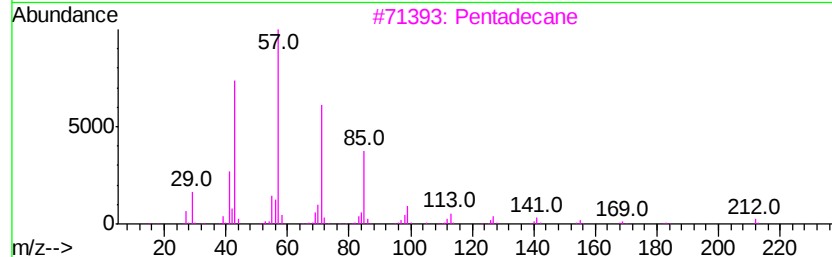
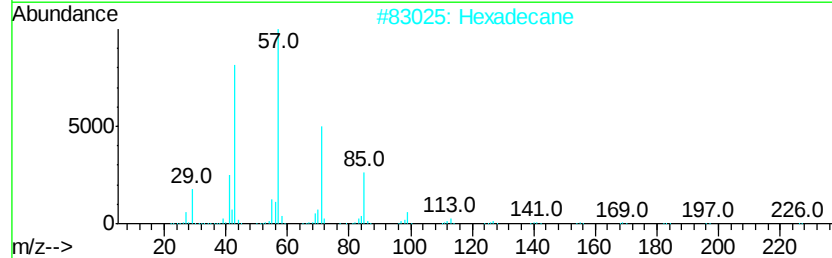
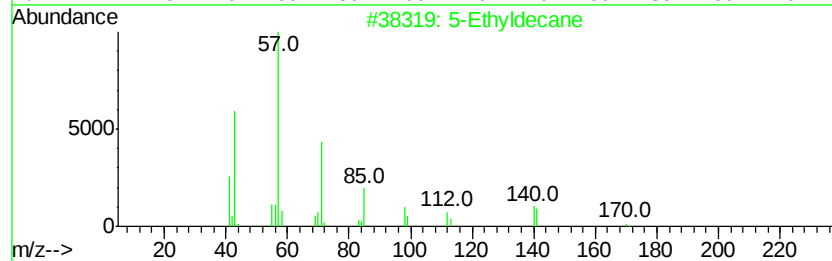
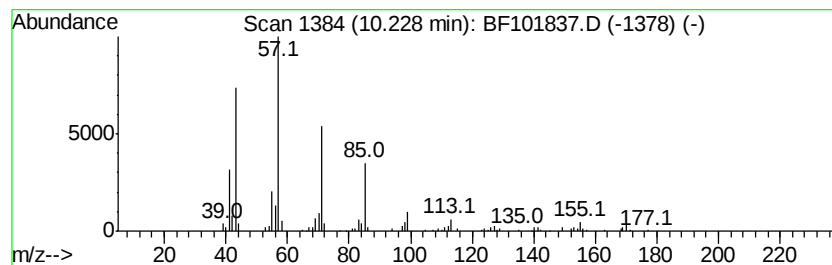
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Peak Number 6 5-Ethyldecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.		
10.23	5.15 ng	214908	Acenaphthene-d10		9.82		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Ethyldecane	170	C12H26	017302-36-2	91
2			Hexadecane	226	C16H34	000544-76-3	90
3			Pentadecane	212	C15H32	000629-62-9	86
4			Tetradecane	198	C14H30	000629-59-4	80
5			Heptadecane	240	C17H36	000629-78-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122917\
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Acq On : 29 Dec 2017 18:53
Operator : SJ/JU
Sample : I7034-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

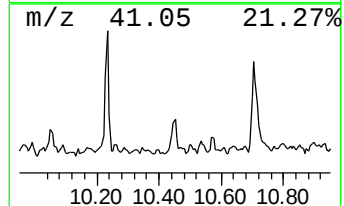
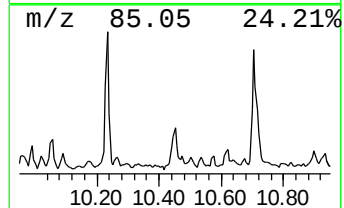
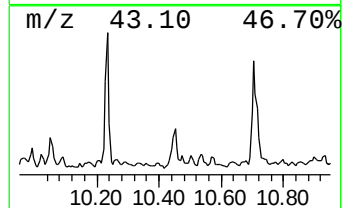
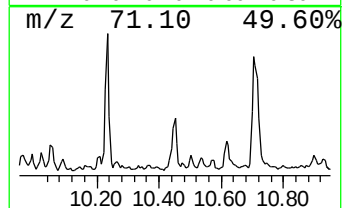
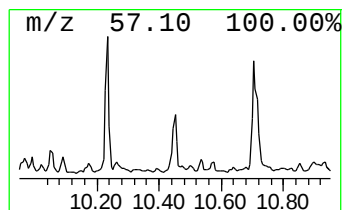
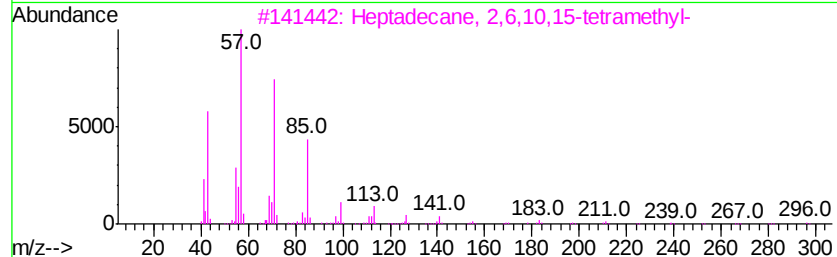
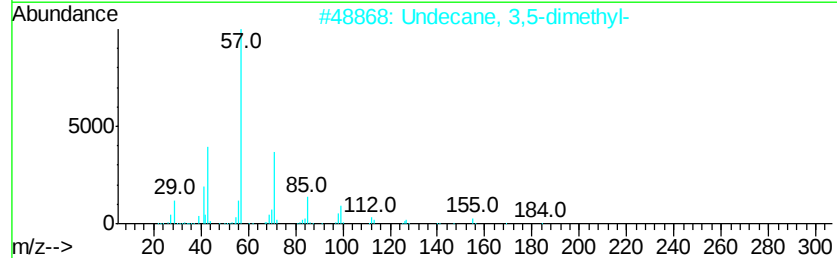
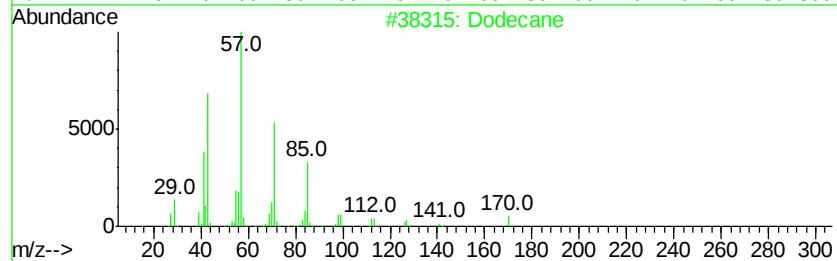
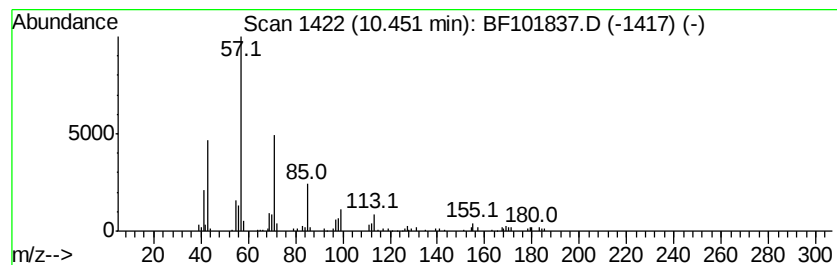
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.45	2.44 ng	101984	Acenaphthene-d10	9.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	83
2	Undecane, 3,5-dimethyl-	184 C13H28	017312-81-1	72
3	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	64
4	1-Octanol, 2-butyl-	186 C12H26O	003913-02-8	59
5	Sulfurous acid, butyl undecyl ester	292 C15H32O3S	1000309-17-8	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122917\
Data File : BF101837.D
Acq On : 29 Dec 2017 18:53
Operator : SJ/JU
Sample : I7034-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TB-05(5-6)

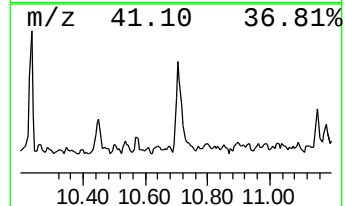
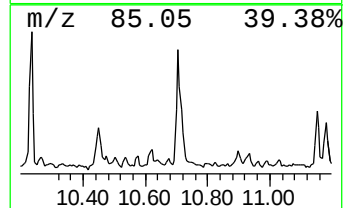
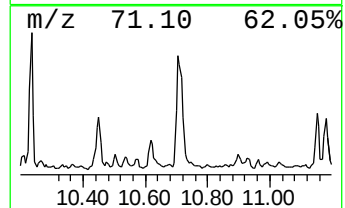
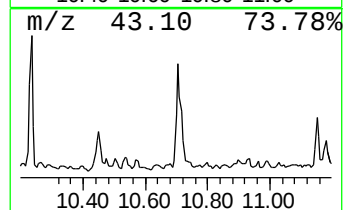
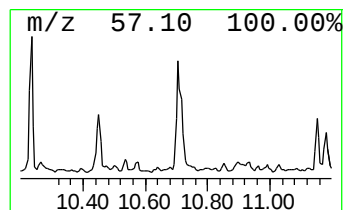
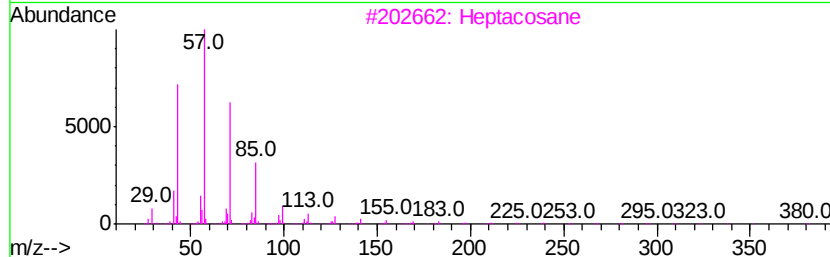
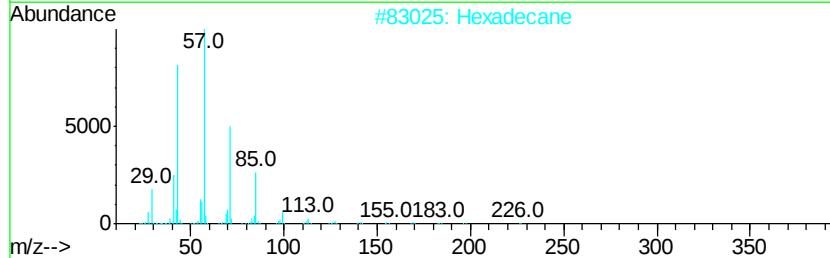
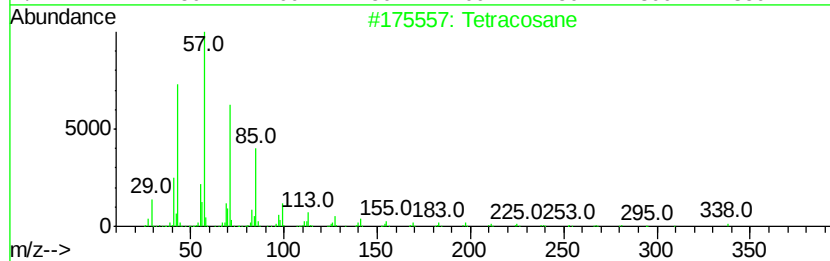
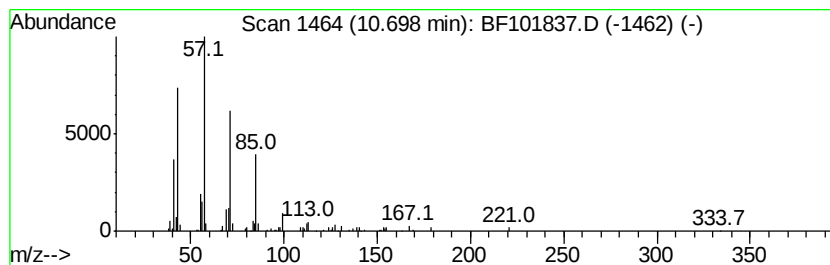
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.70	5.74 ng	225369	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	90
2		Hexadecane	226	C16H34	000544-76-3	90
3		Heptacosane	380	C27H56	000593-49-7	90
4		Tetradecane	198	C14H30	000629-59-4	86
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122917\
Data File : BF101837.D
Acq On : 29 Dec 2017 18:53
Operator : SJ/JU
Sample : I7034-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TB-05(5-6)

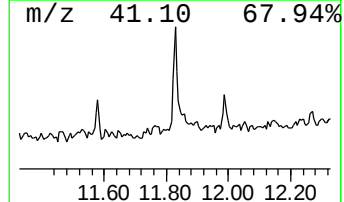
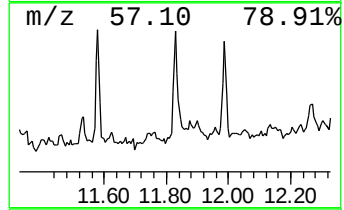
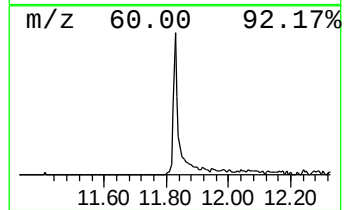
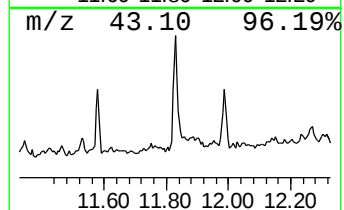
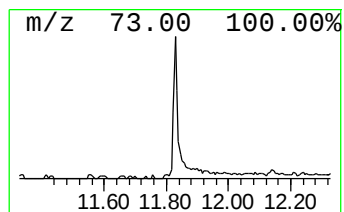
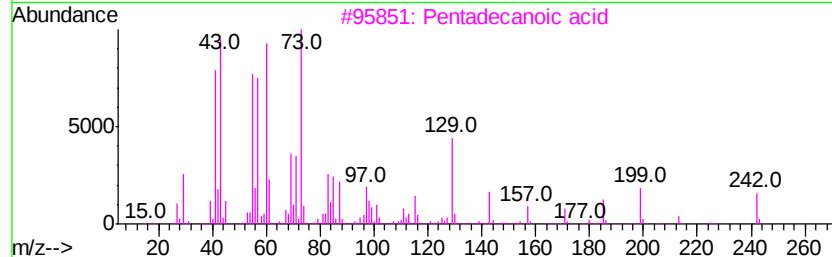
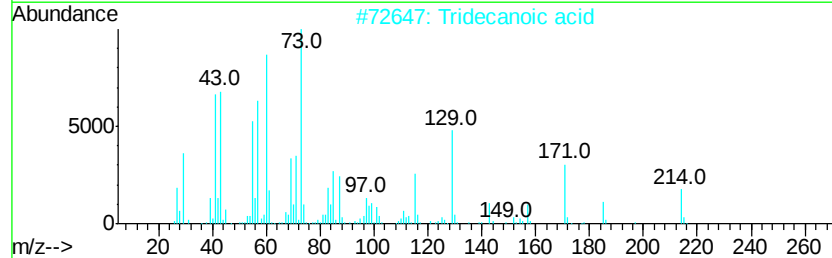
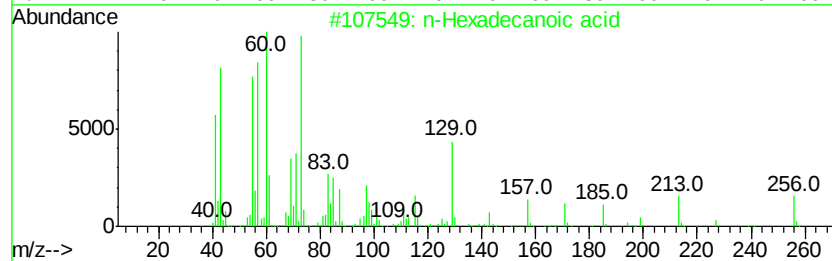
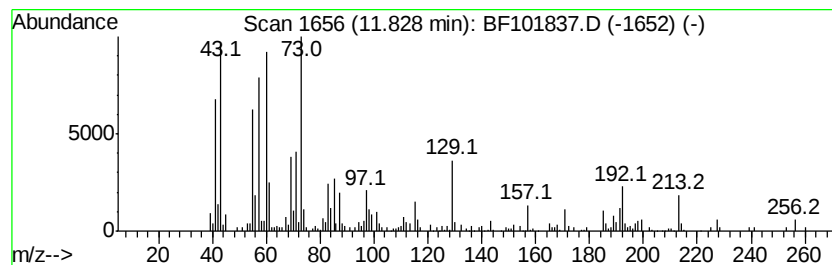
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	5.36 ng	210786	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tridecanoic acid	214	C13H26O2	000638-53-9	95
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	81
5			n-Decanoic acid	172	C10H20O2	000334-48-5	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122917\
Data File : BF101837.D
Acq On : 29 Dec 2017 18:53
Operator : SJ/JU
Sample : I7034-08
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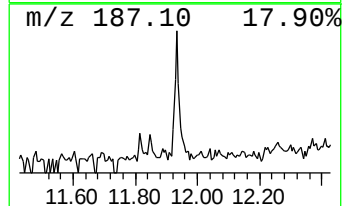
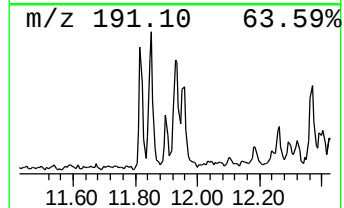
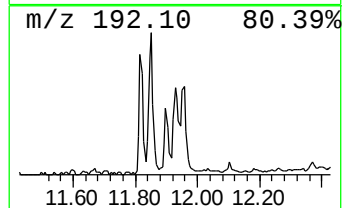
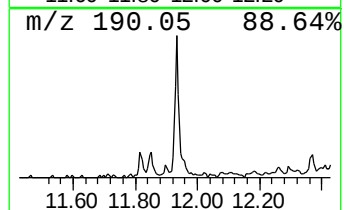
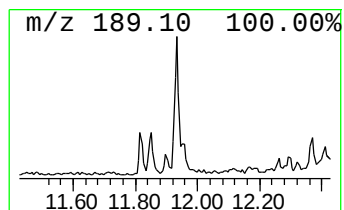
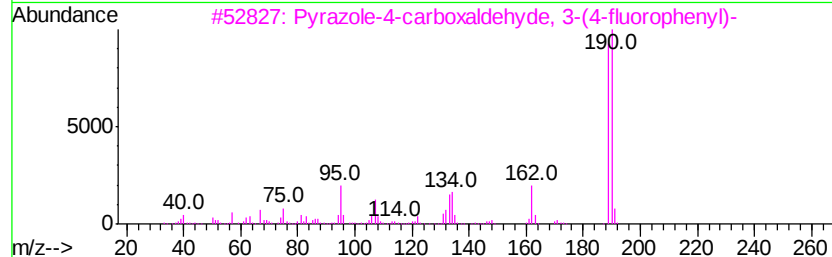
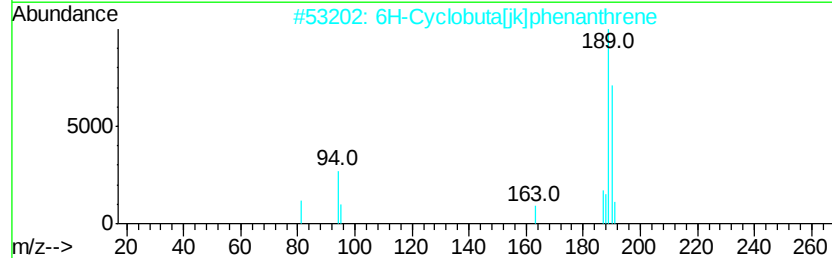
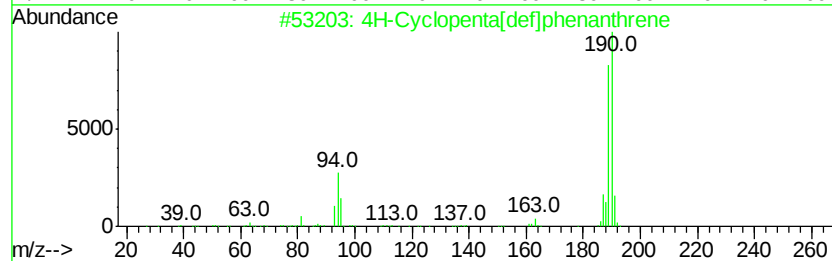
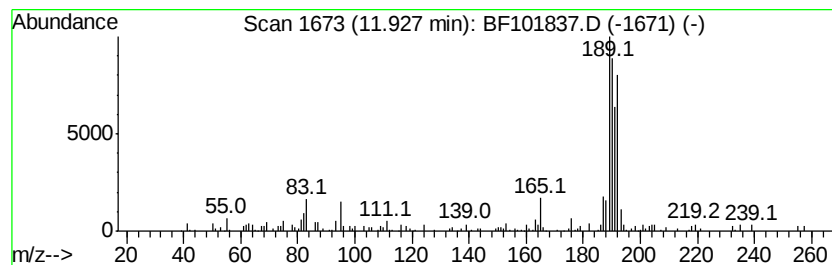
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ClientSampleId :
TB-05(5-6)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.93	2.07 ng	81394	Phenanthrene-d10	11.32		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	58
3		Pvrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	43
4		Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	40
5		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	33



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122917\
Data File : BF101837.D
Acq On : 29 Dec 2017 18:53
Operator : SJ/JU
Sample : I7034-08
Misc :
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Instrument :
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ClientSampled :
TB-05(5-6)

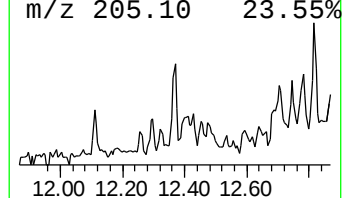
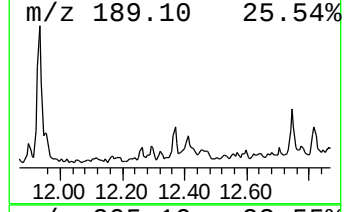
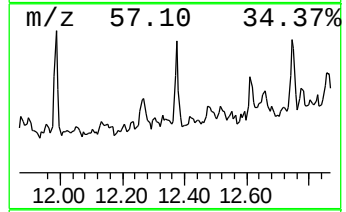
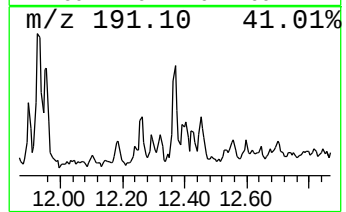
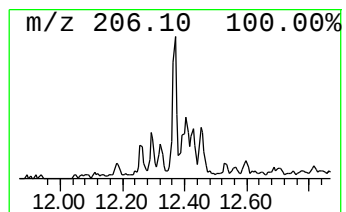
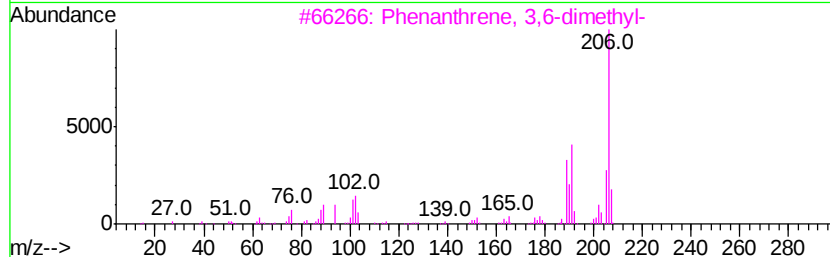
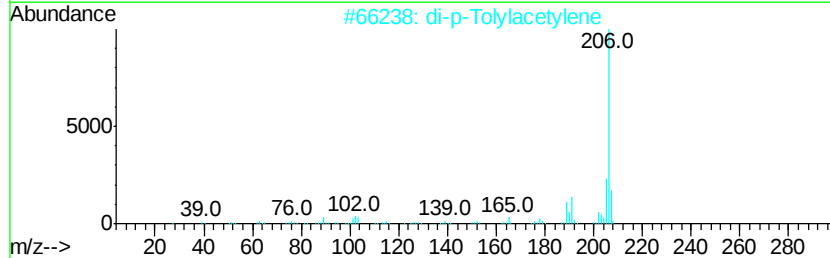
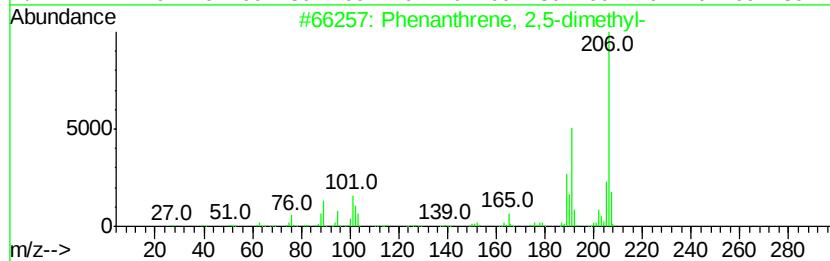
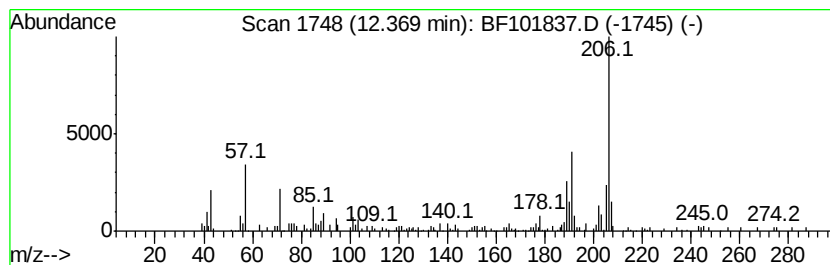
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	2.91 ng	114226	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2		di-p-Tolylacetylene	206	C16H14	002789-88-0	94
3		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
4		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
5		9,10-Dimethylantracene	206	C16H14	000781-43-1	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122917\
Data File : BF101837.D
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ClientSampleId :
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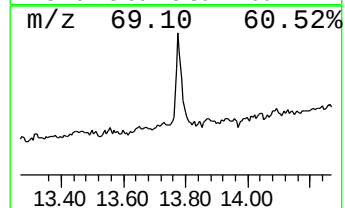
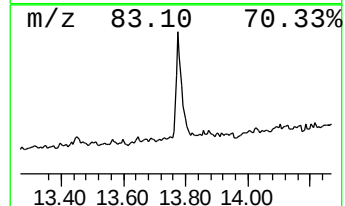
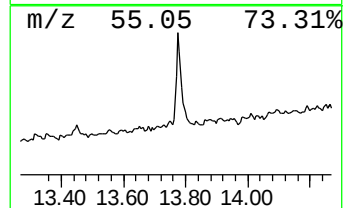
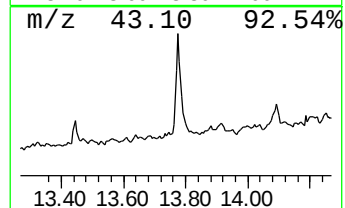
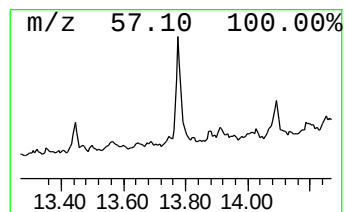
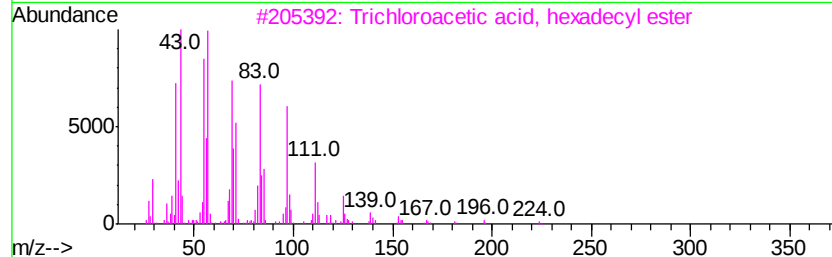
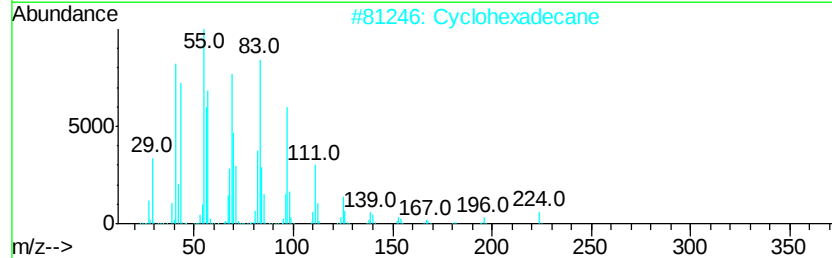
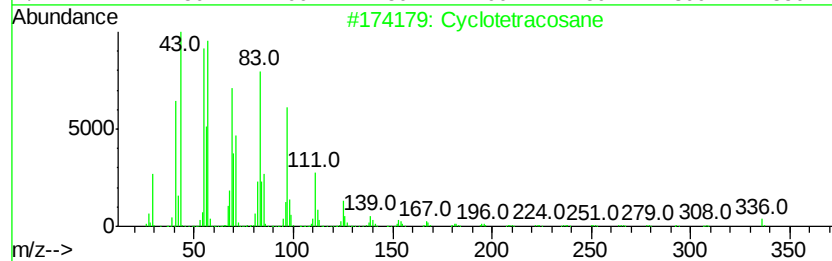
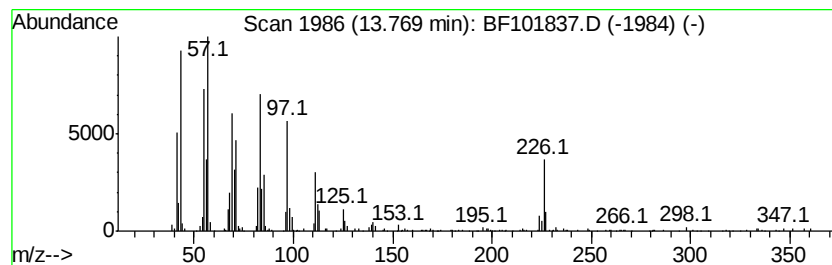
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclotetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	10.12 ng	507072	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	93
2		Cyclohexadecane	224	C16H32	000295-65-8	91
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	89
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	76
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122917\
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 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TB-05(5-6)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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...	5.01	5.8 ng		240912	1	6.79	825474	20.0
unknown6.56	6.56	71.5 ng		2951420	1	6.79	825474	20.0
Tetradecane	9.20	2.6 ng		107251	3	9.82	835209	20.0
Hexadecane	9.73	3.9 ng		163524	3	9.82	835209	20.0
5-Ethyldecane	10.23	5.2 ng		214908	3	9.82	835209	20.0
Dodecane	10.45	2.4 ng		101984	3	9.82	835209	20.0
Tetracosane	10.70	5.7 ng		225369	4	11.32	785831	20.0
n-Hexadecanoic acid	11.83	5.4 ng		210786	4	11.32	785831	20.0
4H-Cyclopenta[def...	11.93	2.1 ng		81394	4	11.32	785831	20.0
Phenanthrene, 2,5...	12.37	2.9 ng		114226	4	11.32	785831	20.0
Cyclotetracosane	13.77	10.1 ng		507072	5	13.97	1002390	20.0