

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
 Data File : BF102849.D
 Acq On : 8 Feb 2018 12:09
 Operator : SJ/JU
 Sample : J1469-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WABUT-3

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-BF020318.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	2.163	8	13	22	rVB	459602	589844	11.71%	1.173%
2	5.116	512	515	524	rVB	157564	201655	4.00%	0.401%
3	5.504	575	581	584	rBV	4848541	5037886	100.00%	10.014%
4	6.510	746	752	755	rBV	4509149	5005292	99.35%	9.950%
5	6.651	771	776	780	rBV	4724139	4921616	97.69%	9.783%
6	6.881	812	815	819	rBV	1660716	1375904	27.31%	2.735%
7	7.040	838	842	845	rBV	4048338	3705514	73.55%	7.366%
8	7.445	906	911	914	rBV	3422661	3373447	66.96%	6.706%
9	8.163	1029	1033	1037	rBV	2065830	1833074	36.39%	3.644%
10	9.239	1211	1216	1219	rBV	4884520	4743735	94.16%	9.430%
11	9.316	1226	1229	1231	rBV	140108	113169	2.25%	0.225%
12	9.628	1279	1282	1285	rBV	494827	474955	9.43%	0.944%
13	9.781	1305	1308	1311	rBV2	54021	59415	1.18%	0.118%
14	9.845	1311	1319	1321	rBV	369156	330831	6.57%	0.658%
15	9.922	1328	1332	1336	rBV	2175754	1850157	36.72%	3.678%
16	10.075	1354	1358	1361	rBV2	44575	60435	1.20%	0.120%
17	10.163	1371	1373	1378	rBV3	81725	80500	1.60%	0.160%
18	10.345	1401	1404	1407	rVB	471832	363500	7.22%	0.723%
19	10.563	1436	1441	1447	rBV	210872	266159	5.28%	0.529%
20	10.645	1453	1455	1460	rBV2	50162	59549	1.18%	0.118%
21	10.710	1462	1466	1472	rVB	2735167	2761556	54.82%	5.489%
22	10.816	1480	1484	1490	rBV2	750318	937868	18.62%	1.864%
23	11.010	1514	1517	1519	rBV2	48004	52631	1.04%	0.105%
24	11.263	1557	1560	1563	rBV	429702	377843	7.50%	0.751%
25	11.292	1563	1565	1572	rVB	264311	267329	5.31%	0.531%
26	11.410	1581	1585	1588	rBV	2012407	1778655	35.31%	3.536%
27	11.433	1588	1589	1595	rVB4	49927	70021	1.39%	0.139%
28	11.645	1622	1625	1628	rVB2	61742	54605	1.08%	0.109%
29	11.692	1630	1633	1636	rVB	359099	267569	5.31%	0.532%
30	11.927	1669	1673	1681	rBV	421284	437064	8.68%	0.869%
31	12.098	1699	1702	1705	rBV	209609	160189	3.18%	0.318%
32	12.251	1726	1728	1737	rVB2	53570	53976	1.07%	0.107%
33	12.486	1766	1768	1772	rVB	90826	73575	1.46%	0.146%
34	12.622	1787	1791	1796	rBV	177394	164144	3.26%	0.326%

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

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

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35	12.710	1803	1806	1810	rBV3	57318	69558	1.38%	0.138%
36	12.851	1828	1830	1833	rVB	116325	100340	1.99%	0.199%
37	12.880	1833	1835	1842	rVB3	73728	83599	1.66%	0.166%
38	12.992	1850	1854	1858	rBV	3372789	3501596	69.51%	6.961%
39	13.545	1945	1948	1952	rBV3	34200	56631	1.12%	0.113%
40	13.874	2001	2004	2012	rBV	811705	829370	16.46%	1.649%
41	14.045	2029	2033	2036	rBV	1157462	1215670	24.13%	2.417%
42	14.069	2036	2037	2043	rVB	145697	133096	2.64%	0.265%
43	14.198	2057	2059	2067	rVB5	39446	55075	1.09%	0.109%
44	14.510	2109	2112	2117	rBV3	110768	125291	2.49%	0.249%
45	15.086	2207	2210	2214	rBV	172070	272744	5.41%	0.542%
46	15.157	2219	2222	2225	rBV2	68930	84220	1.67%	0.167%
47	15.398	2260	2263	2269	rVB	97066	132636	2.63%	0.264%
48	15.527	2280	2285	2295	rBV	811011	1159019	23.01%	2.304%
49	15.992	2359	2364	2368	rBV2	87434	136287	2.71%	0.271%
50	16.039	2369	2372	2380	rVB4	85751	156861	3.11%	0.312%
51	17.615	2635	2640	2650	rVB5	122316	290618	5.77%	0.578%

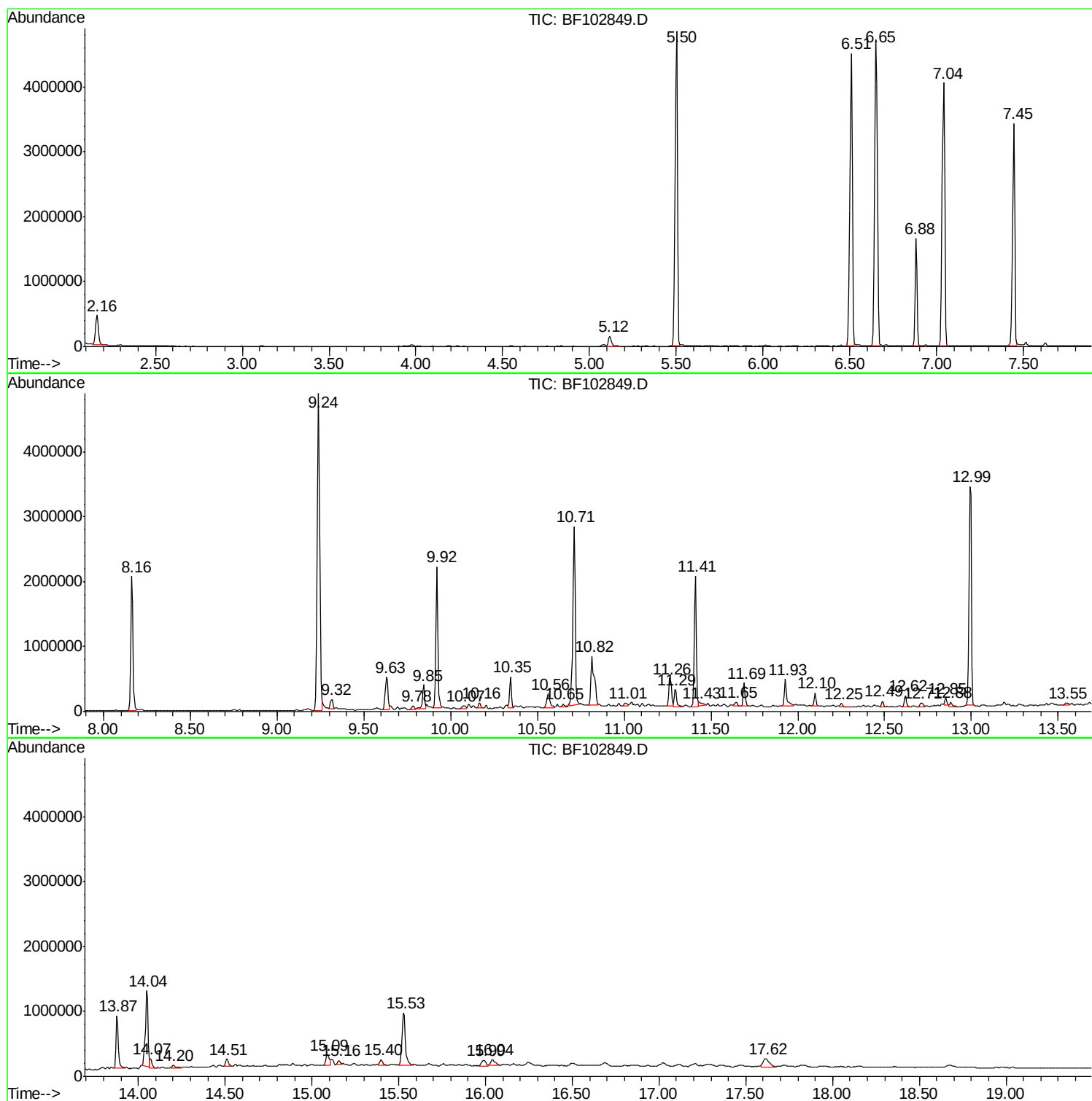
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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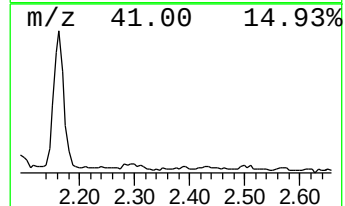
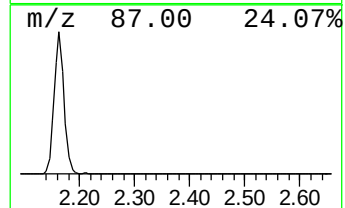
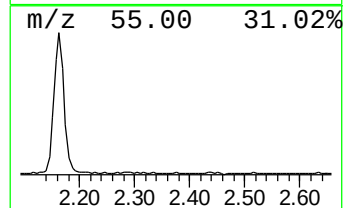
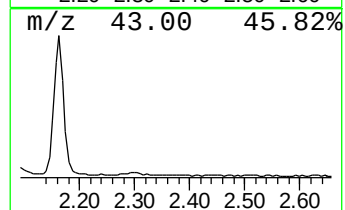
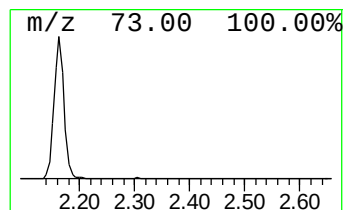
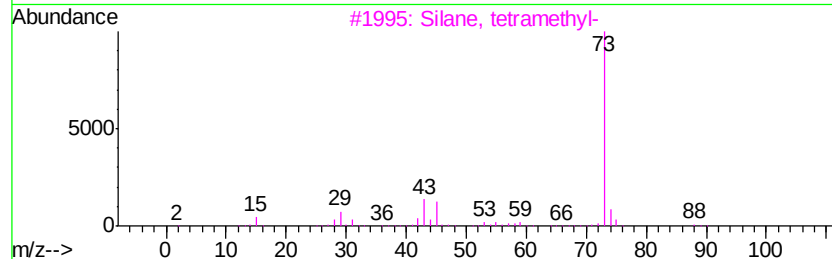
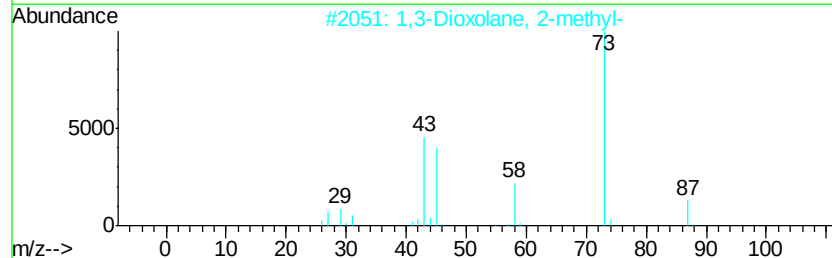
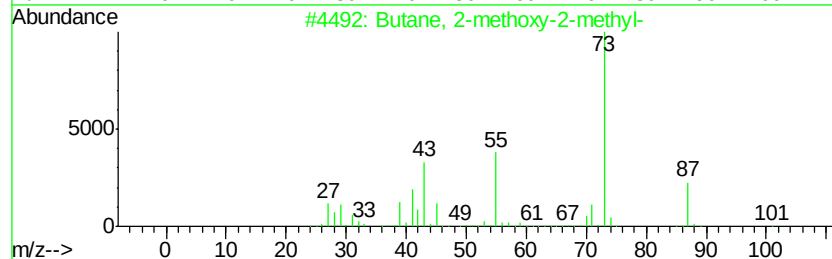
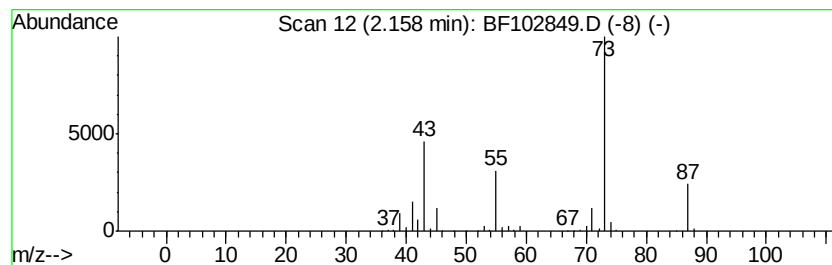
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.16	8.57 ng	589844	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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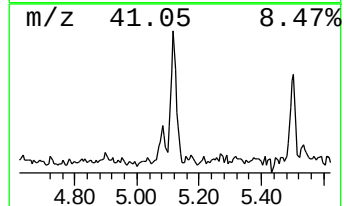
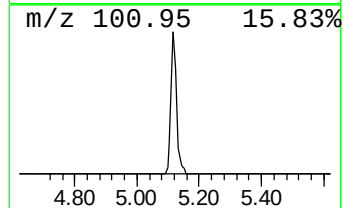
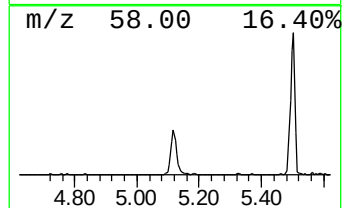
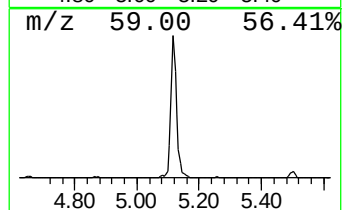
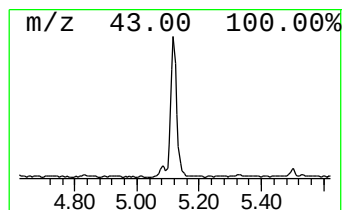
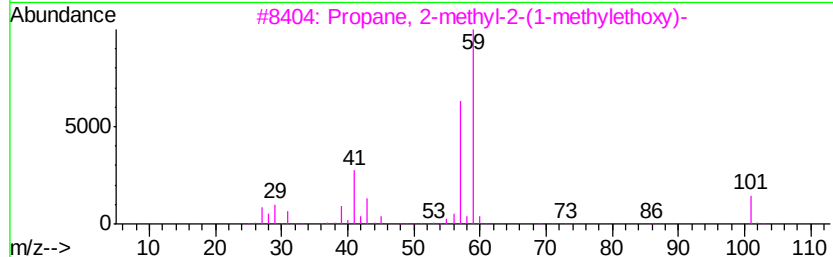
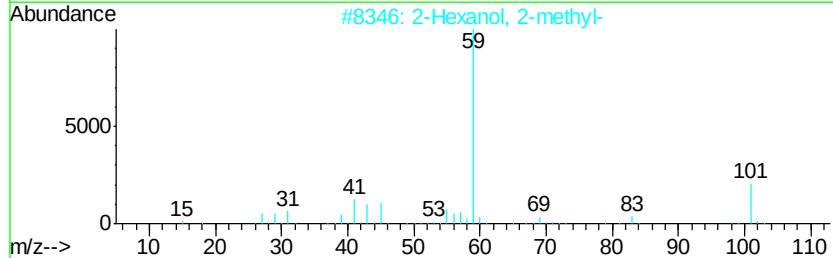
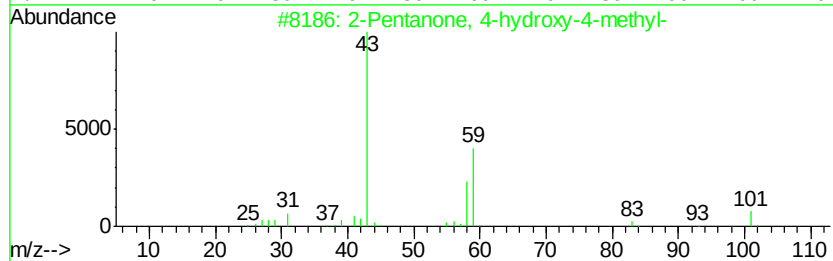
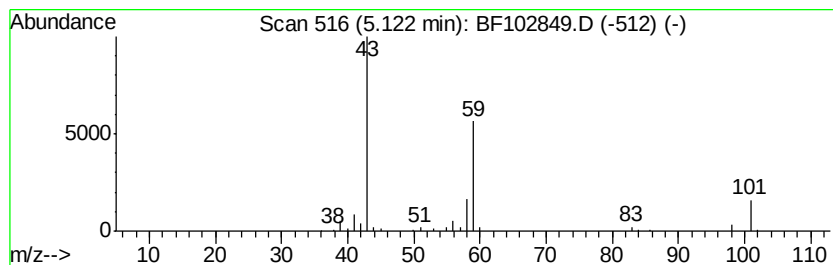
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	2.93 ng	201655	1,4-Dichlorobenzene-d4	6.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	40
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			1,2-Butanediol	90	C4H10O2	000584-03-2	28



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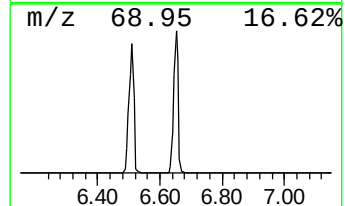
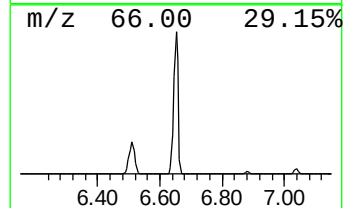
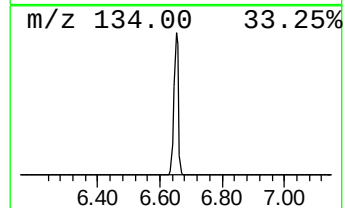
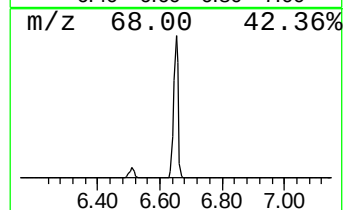
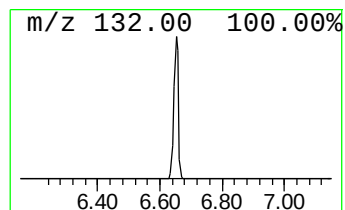
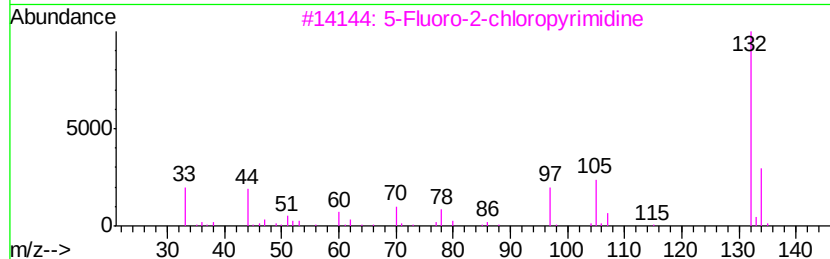
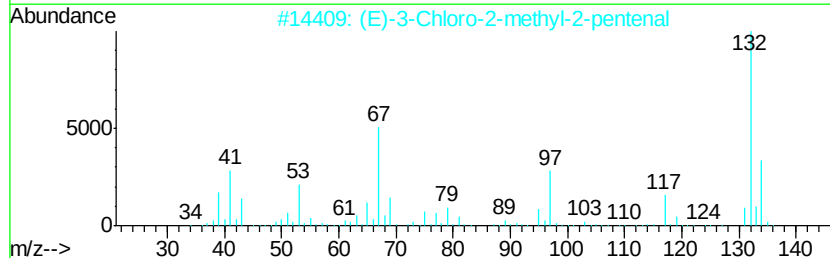
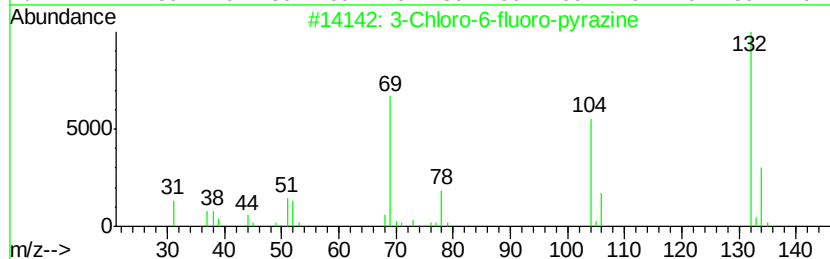
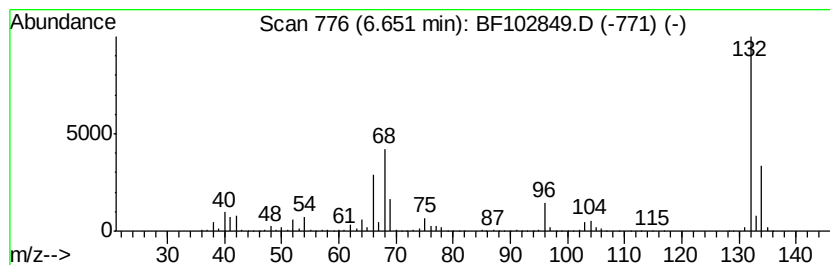
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	71.54 ng	4921620	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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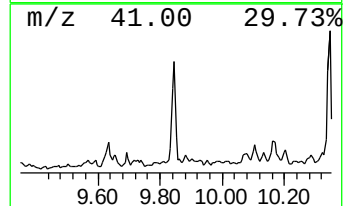
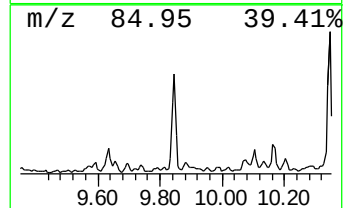
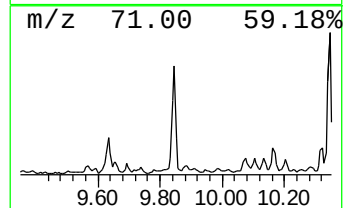
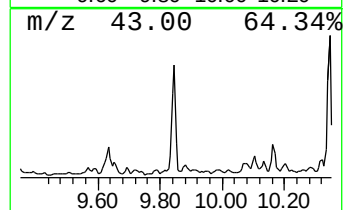
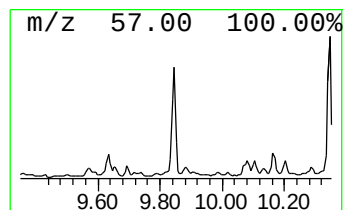
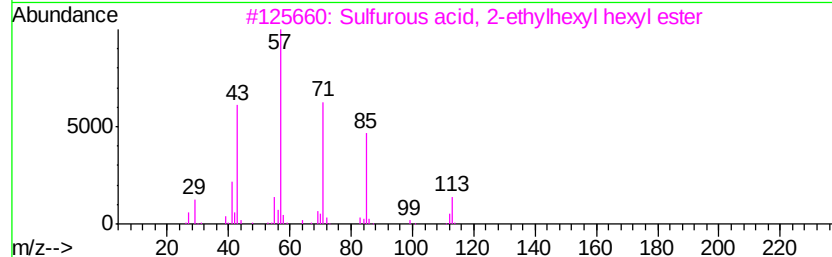
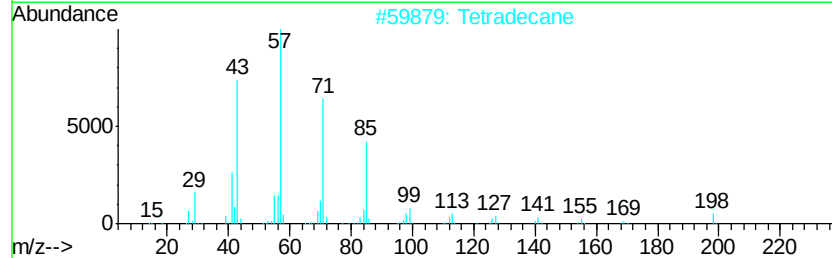
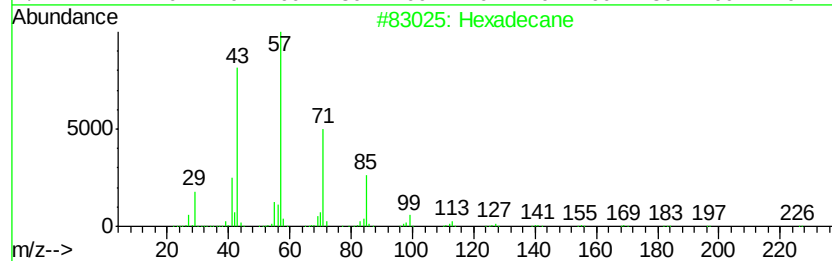
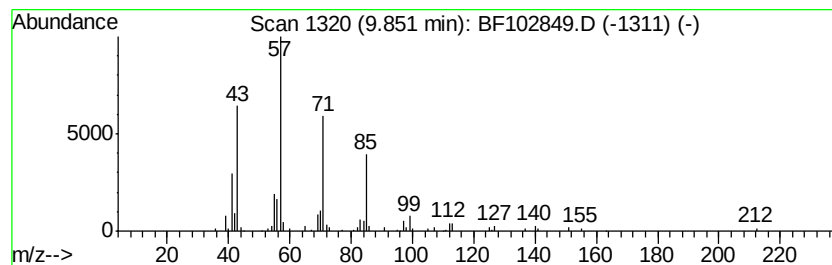
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	3.58 ng	330831	Acenaphthene-d10	9.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	91
2	Tetradecane	198 C14H30	000629-59-4	86
3	Sulfurous acid, 2-ethylhexyl hex...	278 C14H30O3S	1000309-20-2	86
4	2-Bromo dodecane	248 C12H25Br	013187-99-0	72
5	Tetradecane, 4-methyl-	212 C15H32	025117-24-2	72



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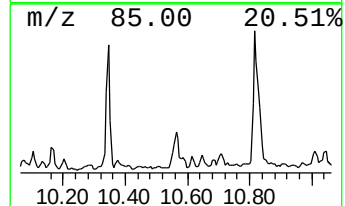
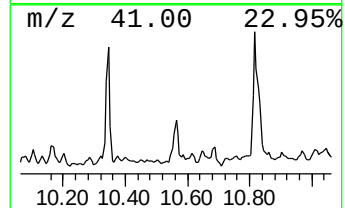
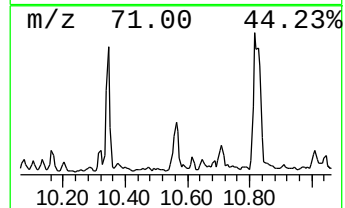
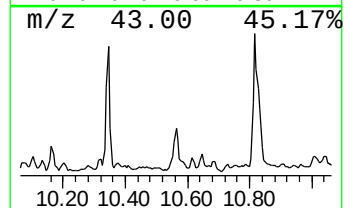
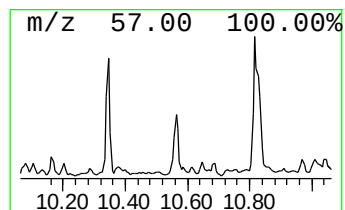
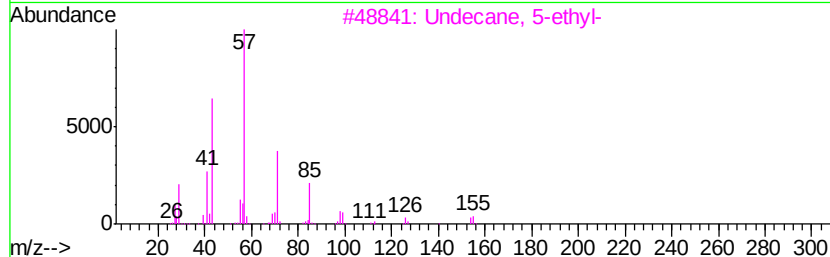
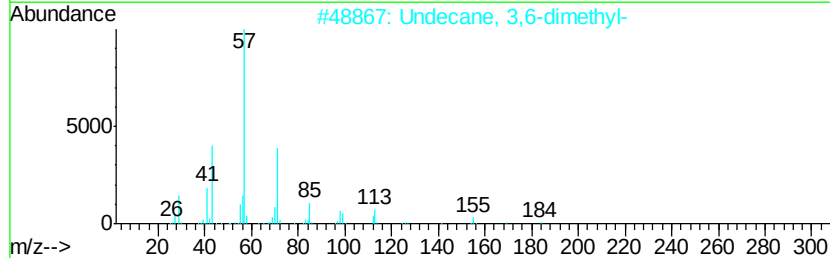
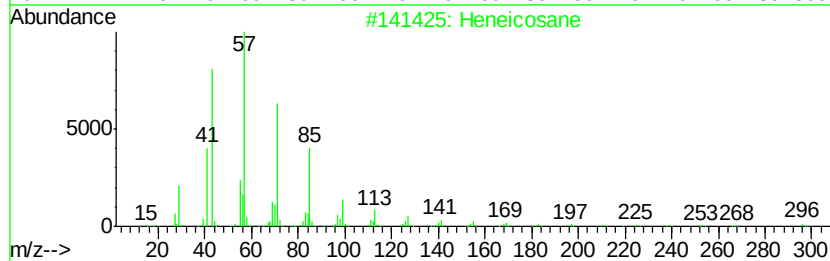
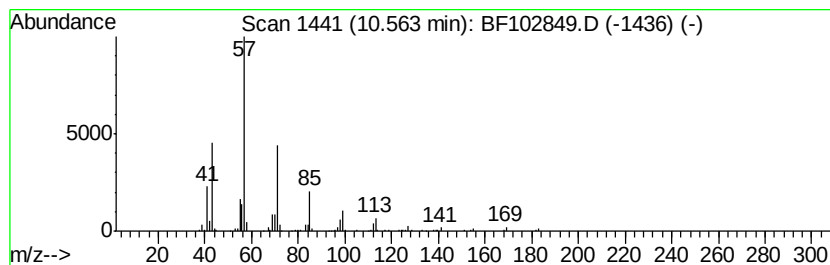
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.56	2.88 ng	266159	Acenaphthene-d10	9.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	80
2	Undecane, 3,6-dimethyl-	184 C13H28	017301-28-9	78
3	Undecane, 5-ethyl-	184 C13H28	017453-94-0	78
4	Octacosane	394 C28H58	000630-02-4	72
5	Decane, 2,6,8-trimethyl-	184 C13H28	062108-26-3	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102849.D
Acq On : 8 Feb 2018 12:09
Operator : SJ/JU
Sample : J1469-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

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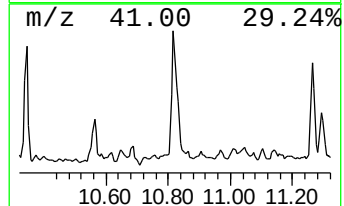
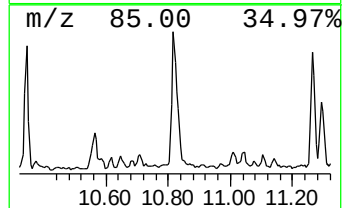
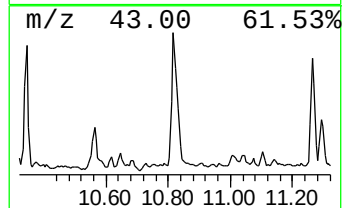
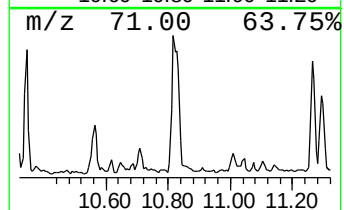
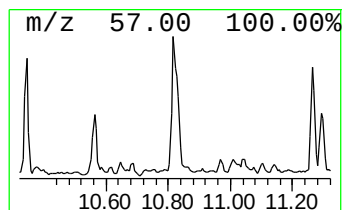
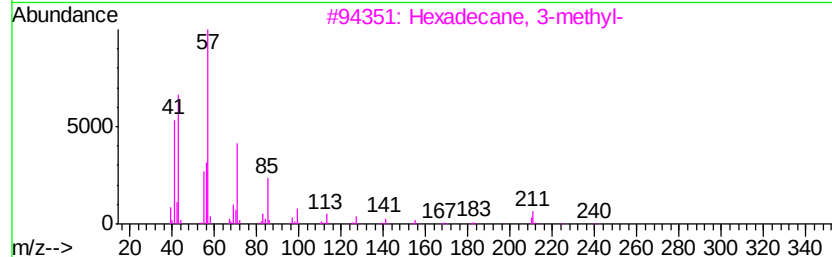
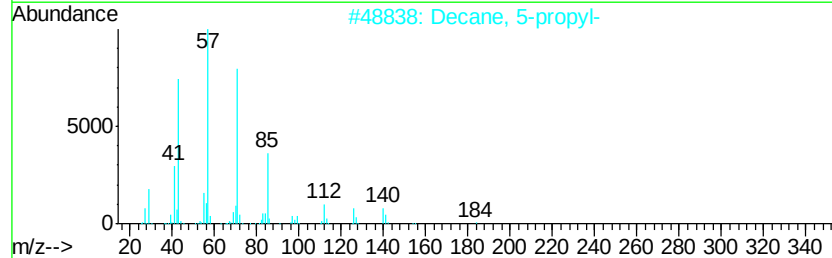
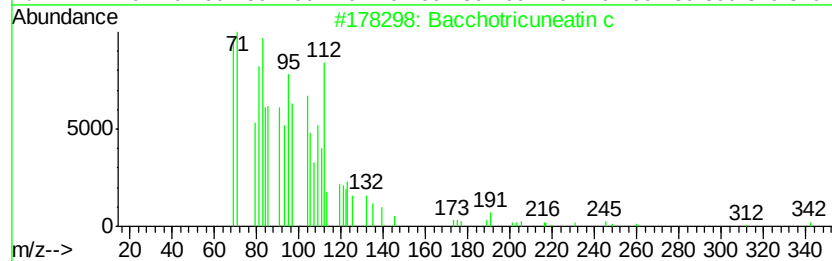
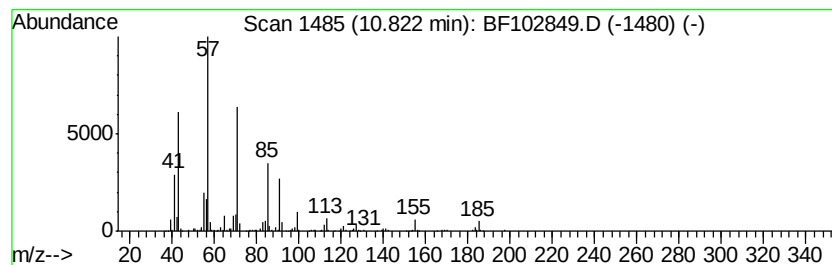
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Bacchotricuneatin c Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.82	10.55 ng	937868	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bacchotricuneatin c	342	C20H22O5	066563-30-2	96
2		Decane, 5-propyl-	184	C13H28	017312-62-8	64
3		Hexadecane, 3-methyl-	240	C17H36	006418-43-5	58
4		Hexadecane	226	C16H34	000544-76-3	52
5		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
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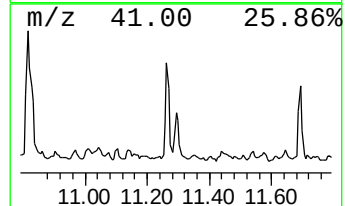
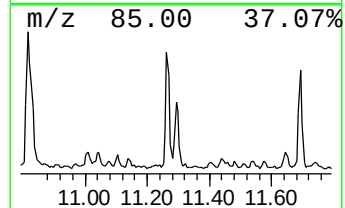
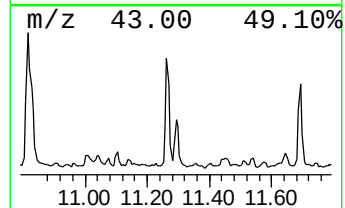
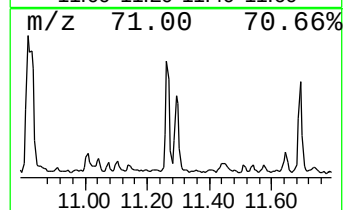
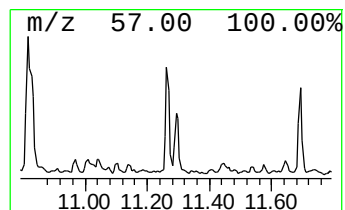
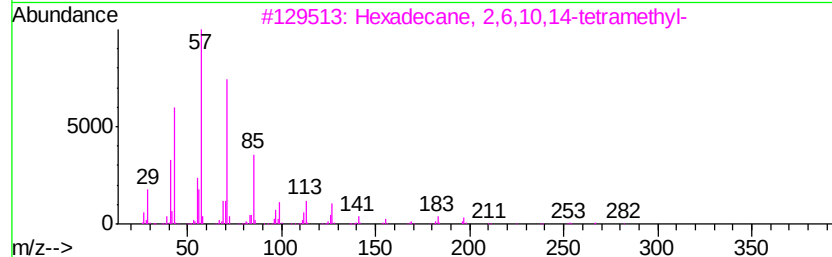
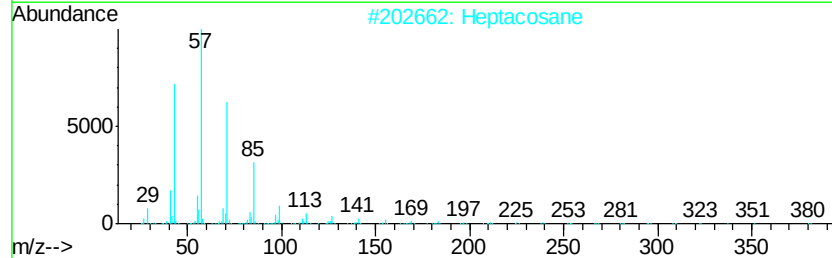
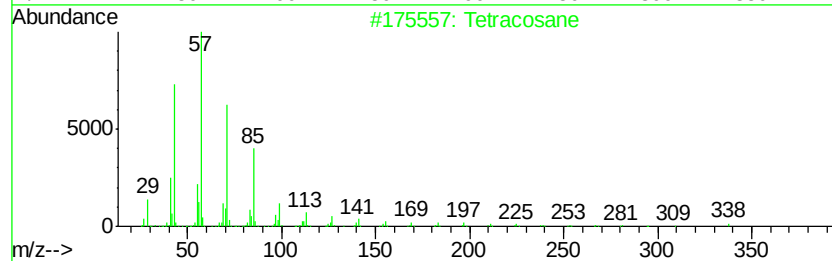
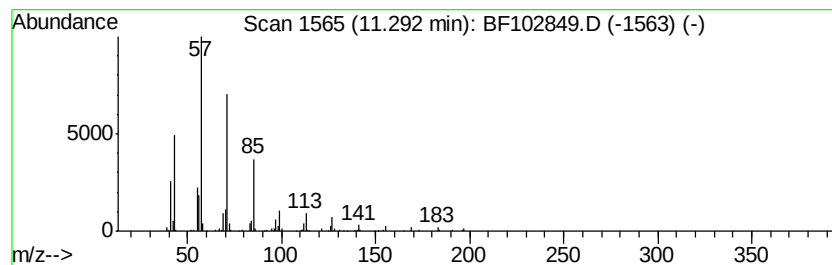
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	3.01 ng	267329	Phenanthrene-d10	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Heptacosane	380	C27H56	000593-49-7	91
3	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
4	Pentacosane	352	C25H52	000629-99-2	90
5	Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
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Operator : SJ/JU
Sample : J1469-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

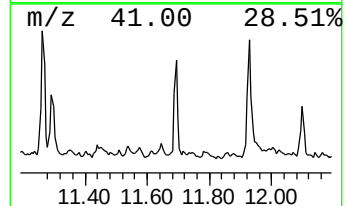
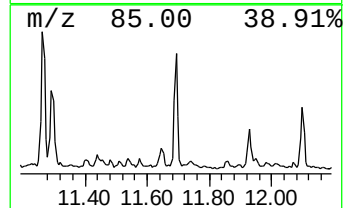
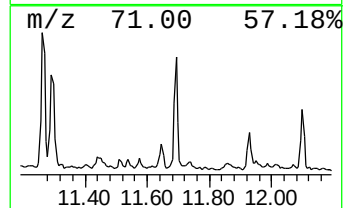
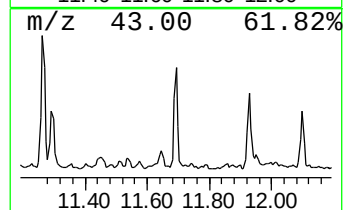
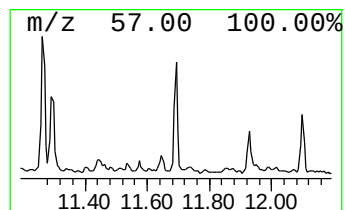
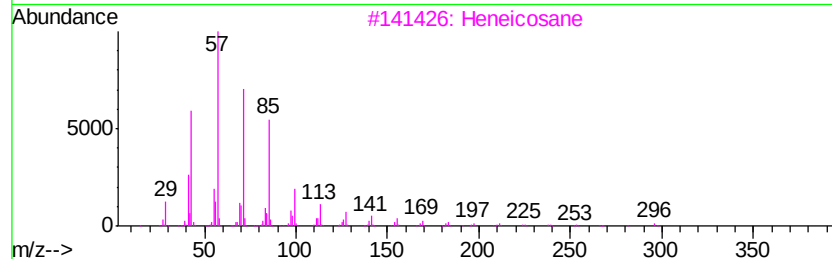
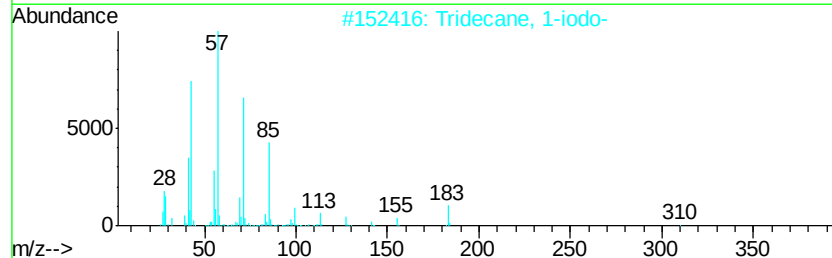
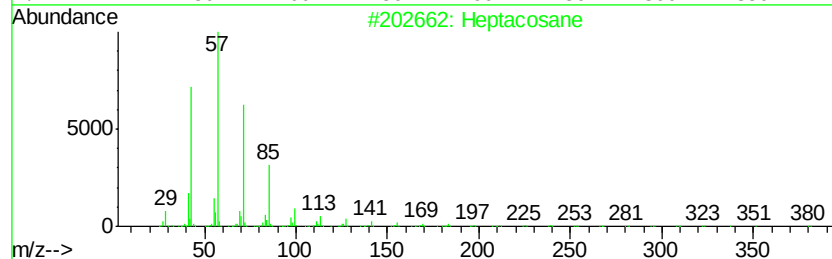
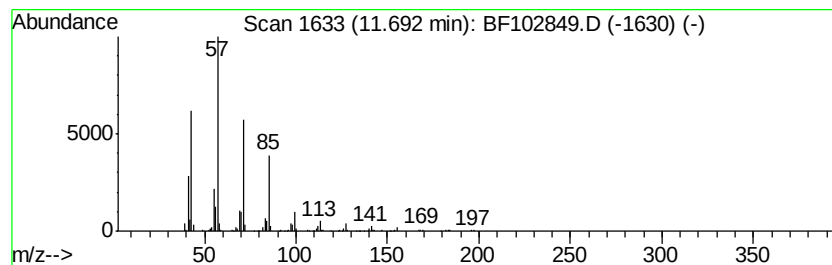
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.69	3.01 ng	267569	Phenanthrene-d10		11.41
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
3	Heneicosane	296	C21H44	000629-94-7	90
4	Pentadecane	212	C15H32	000629-62-9	87
5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
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Acq On : 8 Feb 2018 12:09
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Misc :
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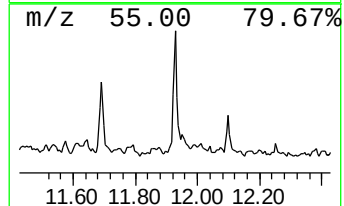
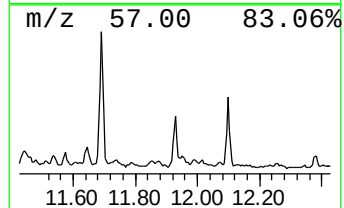
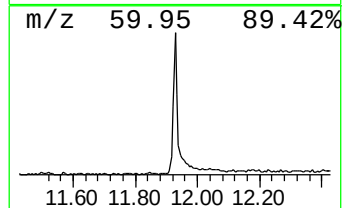
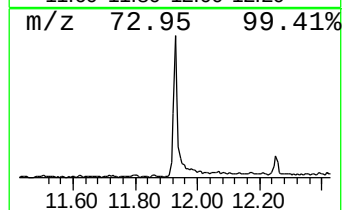
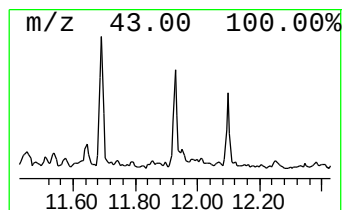
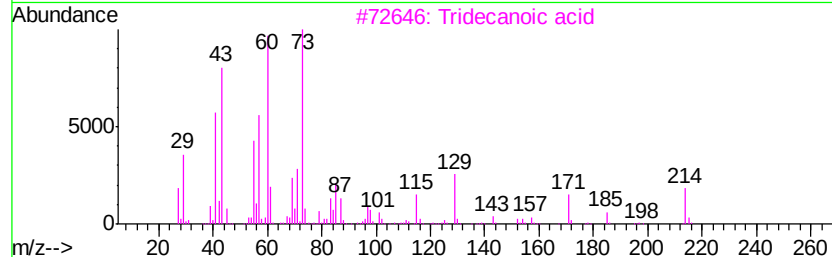
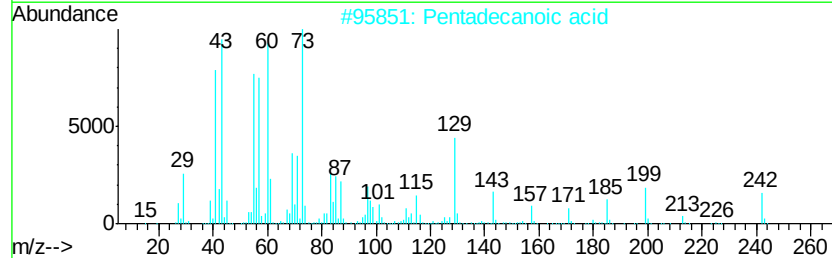
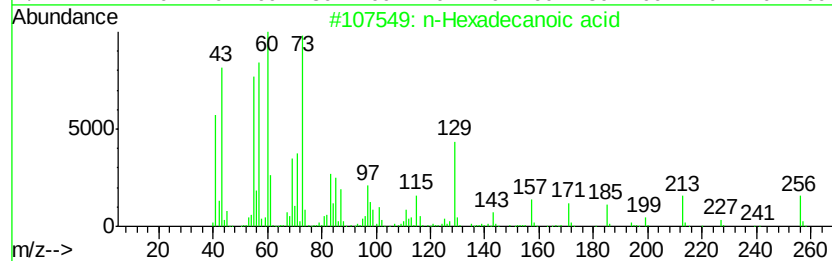
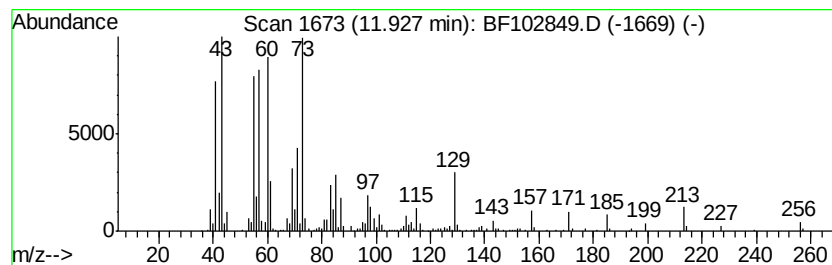
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	4.91 ng	437064	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tridecanoic acid	214	C13H26O2	000638-53-9	81
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		Isoamyl nitrite	117	C5H11NO2	000110-46-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102849.D
Acq On : 8 Feb 2018 12:09
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Misc :
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ClientSampleId :
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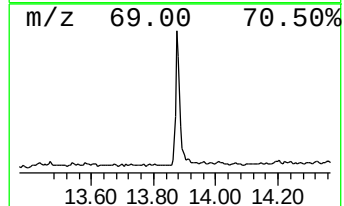
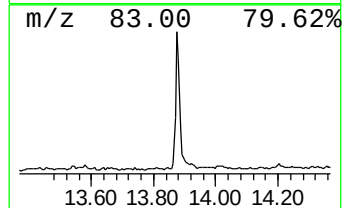
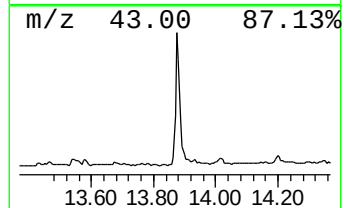
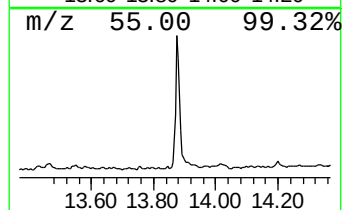
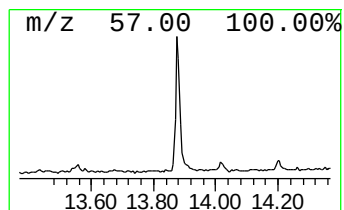
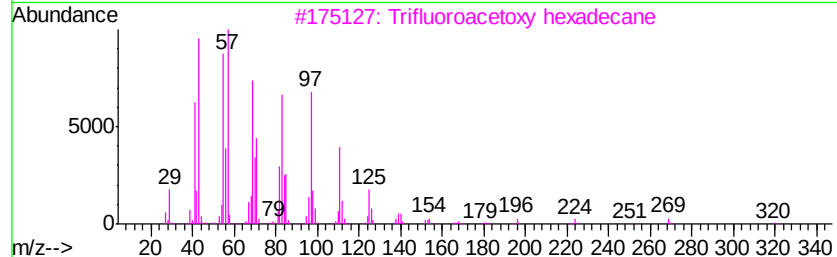
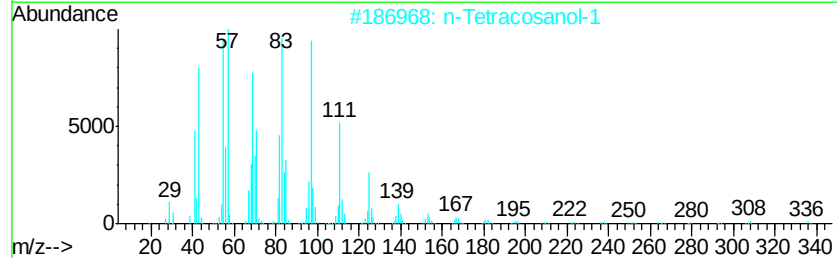
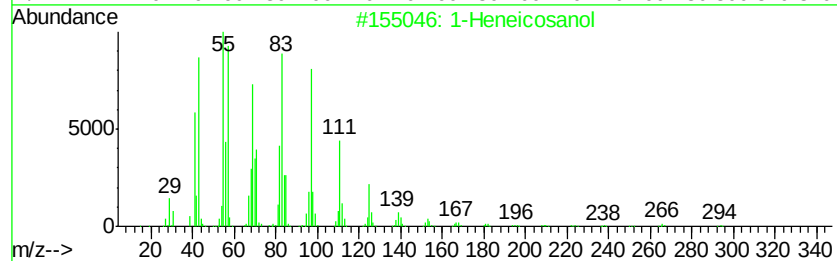
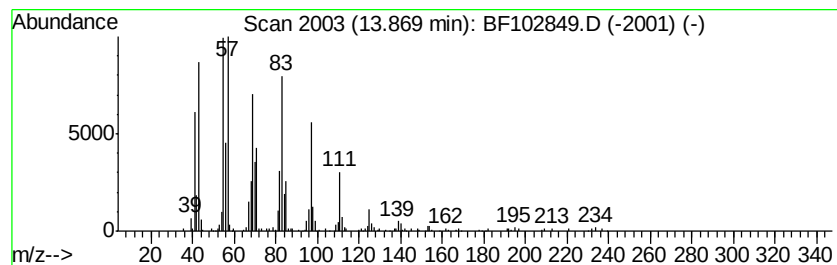
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	13.64 ng	829370	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	94
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	91
3		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	91
5		2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102849.D
Acq On : 8 Feb 2018 12:09
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Misc :
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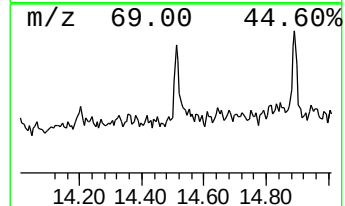
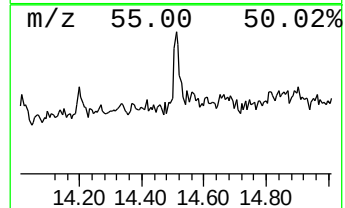
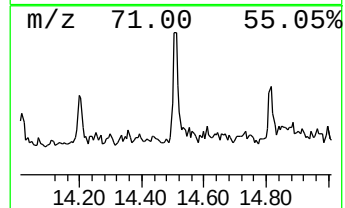
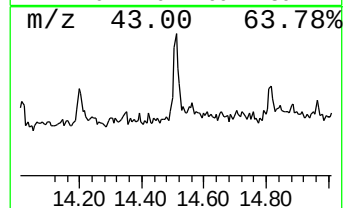
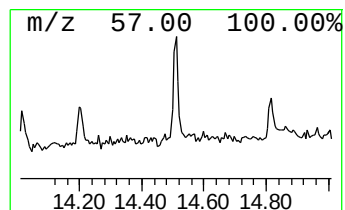
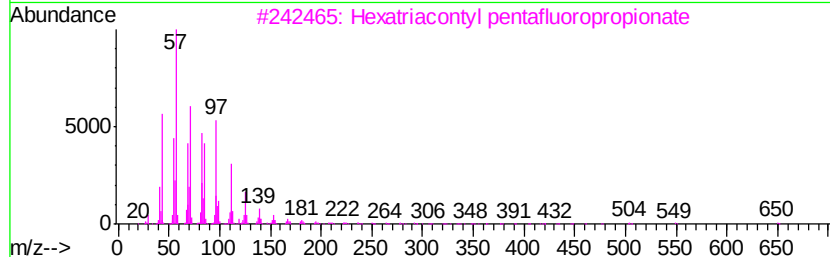
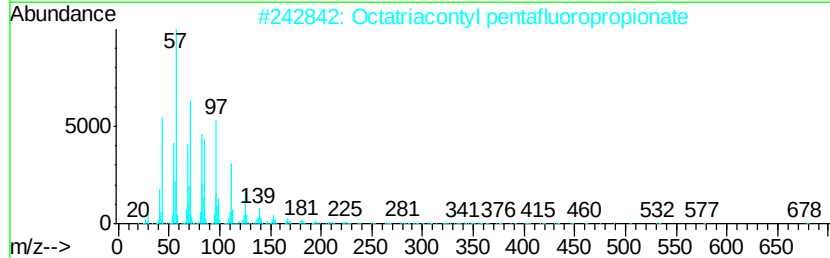
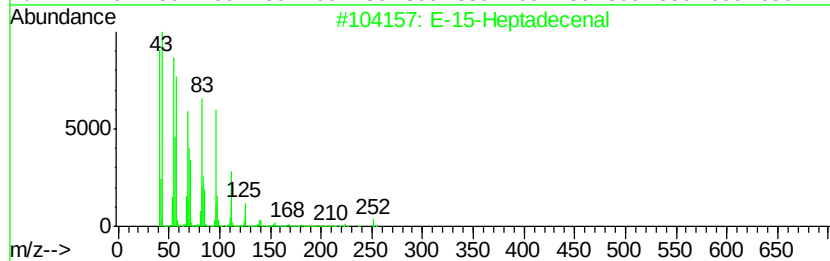
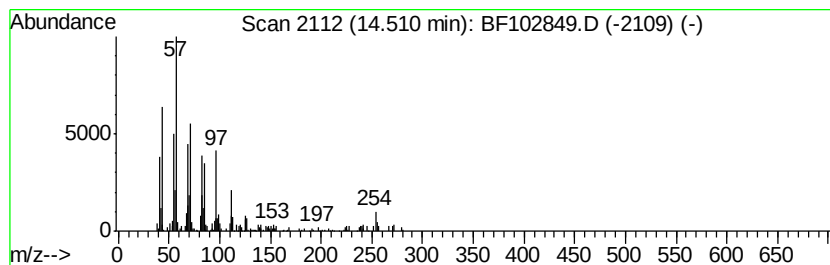
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 E-15-Heptadecenal Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.51	2.06 ng	125291	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	E-15-Heptadecenal	252	C17H32O	1000130-97-9	90
2		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	90
3		Hexatriacontyl pentafluoropropio...	669	C39H73F5O2	1000351-89-0	90
4		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	86
5		Octatriacontyl trifluoroacetate	647	C40H77F3O2	1000351-88-7	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
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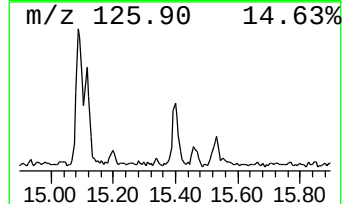
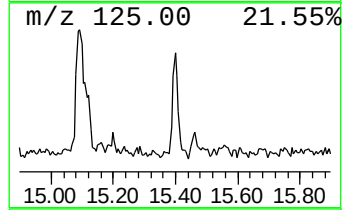
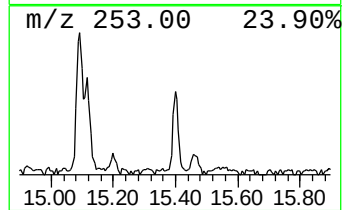
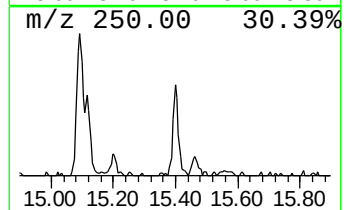
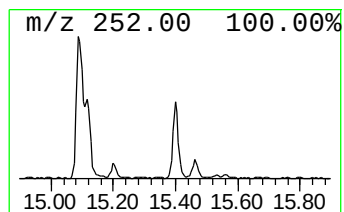
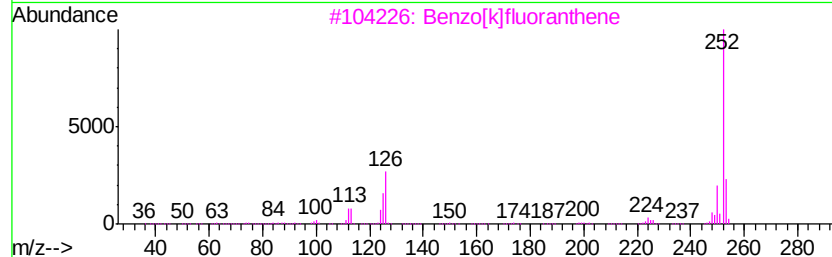
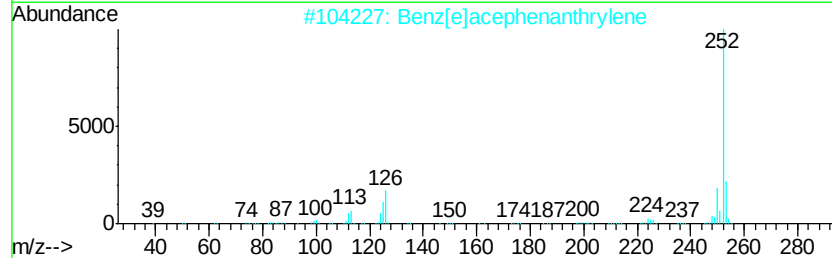
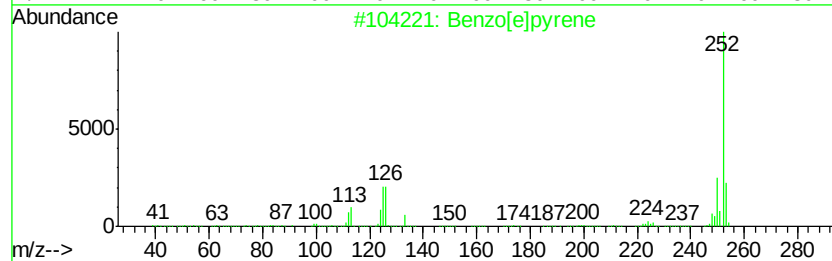
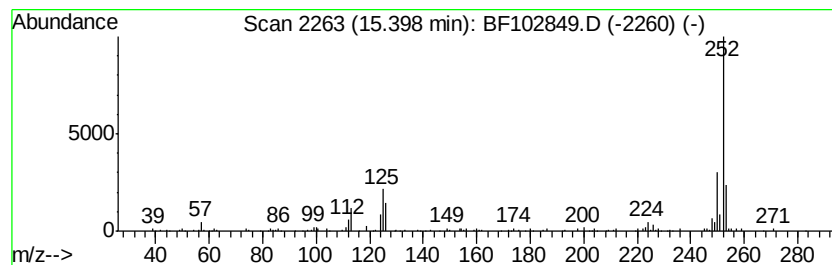
Instrument :
BNA_F
ClientSampleId :
WABUT-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzo[e]pyrene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.40	2.29 ng	132636	Perylene-d12	15.53	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	91
2	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
4	Benzo[j]fluoranthene	252	C20H12	000205-82-3	90
5	Perylene	252	C20H12	000198-55-0	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102849.D
Acq On : 8 Feb 2018 12:09
Operator : SJ/JU
Sample : J1469-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

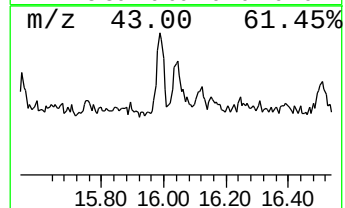
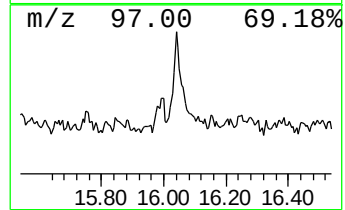
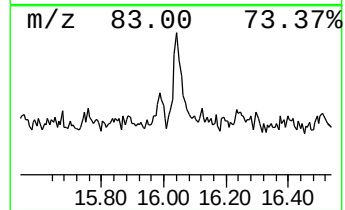
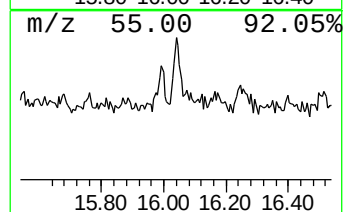
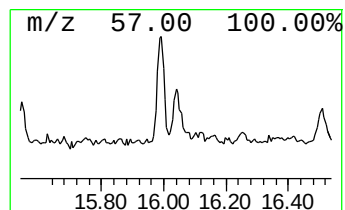
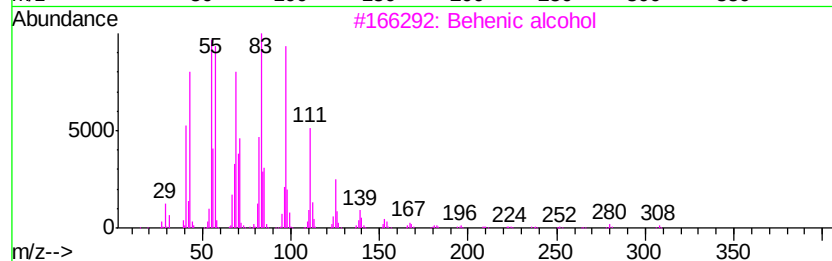
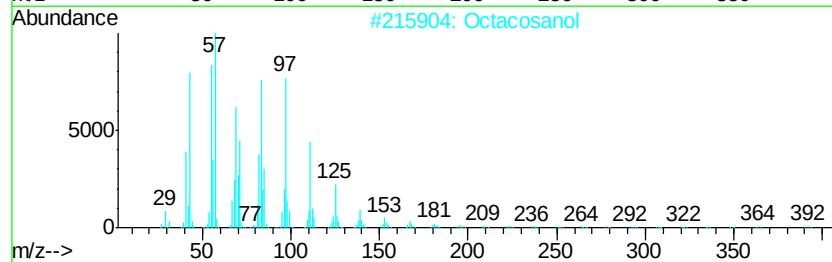
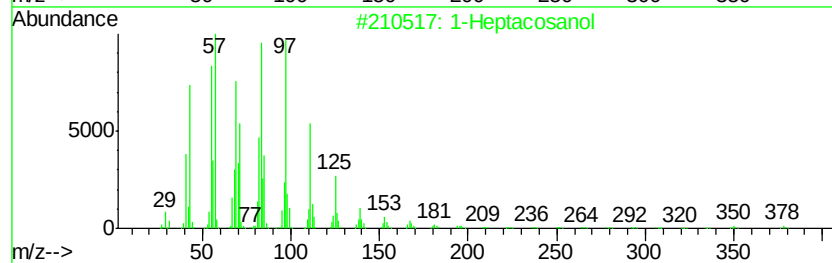
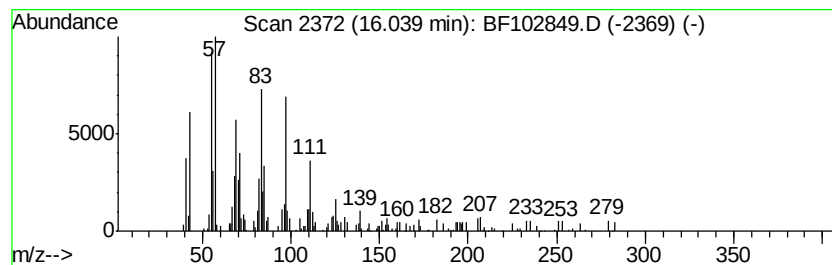
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1-Heptacosanol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.04	2.71 ng	156861	Perylene-d12	15.53	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	90
2	Octacosanol	410	C28H58O	000557-61-9	90
3	Behenic alcohol	326	C22H46O	000661-19-8	87
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	87
5	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
Data File : BF102849.D
Acq On : 8 Feb 2018 12:09
Operator : SJ/JU
Sample : J1469-05
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WABUT-3

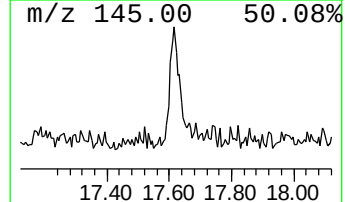
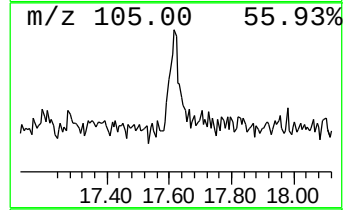
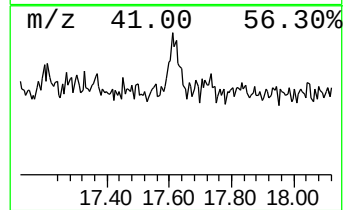
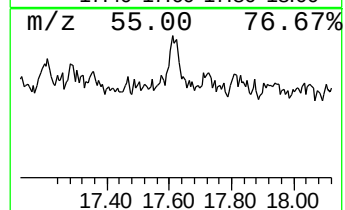
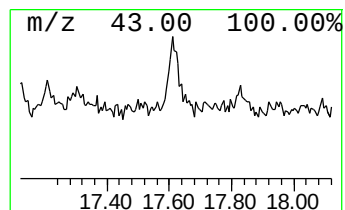
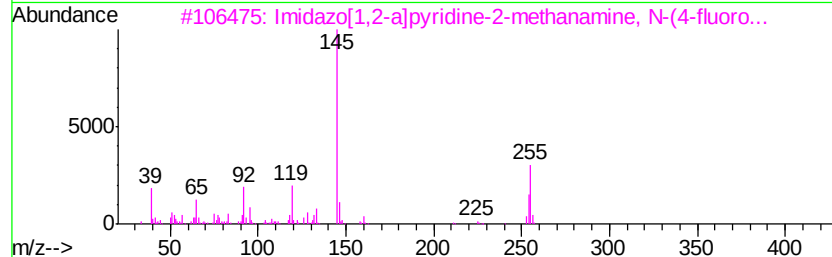
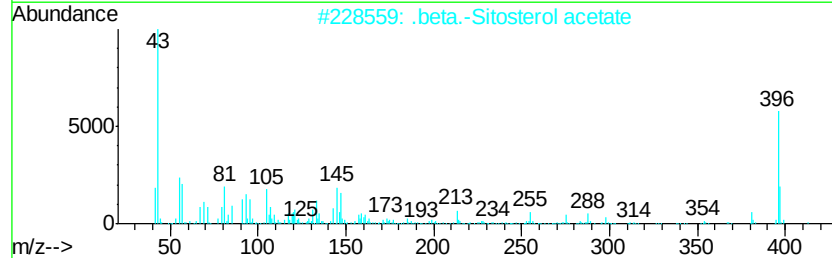
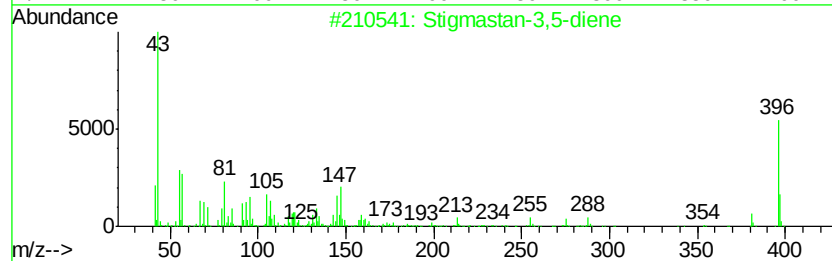
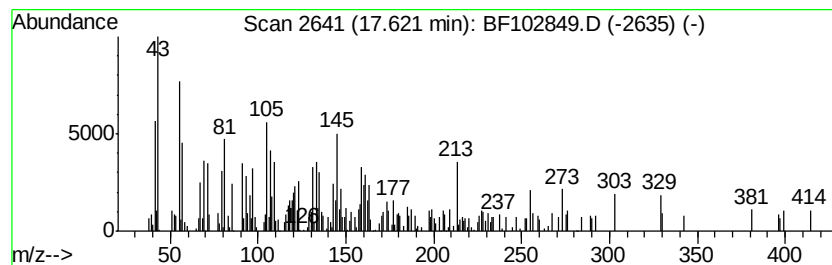
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown17.62 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.62	5.01 ng	290618	Perylene-d12	15.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Stigmastan-3,5-diene	396	C29H48	1000214-16-4	32
2		.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	32
3		Imidazo[1,2-a]pyridine-2-methana...	255	C15H14FN3	1000319-95-6	30
4		Norflurazon	303	C12H9ClF3N3O	027314-13-2	27
5		Benzofuran-2-carboxylic acid, (2...	303	C16H11F2N03	1000311-07-8	16



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020818\
 Data File : BF102849.D
 Acq On : 8 Feb 2018 12:09
 Operator : SJ/JU
 Sample : J1469-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WABUT-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.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...	2.16	8.6	ng	589844	1	6.88	1375900	20.0
2-Pentanone, 4-hy...	5.12	2.9	ng	201655	1	6.88	1375900	20.0
unknown6.65	6.65	71.5	ng	4921620	1	6.88	1375900	20.0
Hexadecane	9.85	3.6	ng	330831	3	9.92	1850160	20.0
Heneicosane	10.56	2.9	ng	266159	3	9.92	1850160	20.0
Bacchotricuneatin c	10.82	10.6	ng	937868	4	11.41	1778660	20.0
Tetracosane	11.29	3.0	ng	267329	4	11.41	1778660	20.0
Heptacosane	11.69	3.0	ng	267569	4	11.41	1778660	20.0
n-Hexadecanoic acid	11.93	4.9	ng	437064	4	11.41	1778660	20.0
1-Heneicosanol	13.87	13.6	ng	829370	5	14.05	1215670	20.0
E-15-Heptadecenal	14.51	2.1	ng	125291	5	14.05	1215670	20.0
Benzo[e]pyrene	15.40	2.3	ng	132636	6	15.53	1159020	20.0
1-Heptacosanol	16.04	2.7	ng	156861	6	15.53	1159020	20.0
unknown17.62	17.62	5.0	ng	290618	6	15.53	1159020	20.0