

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
 Data File : BF111276.D
 Acq On : 11 Dec 2018 14:04
 Operator : JU/SJ
 Sample : J6196-01 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-03-120718-A

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-BF120818.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	4.993	487	494	503	rVB2	224482	307668	5.99%	0.524%
2	5.410	560	565	568	rBV	5170366	4778065	92.98%	8.134%
3	6.428	733	738	750	rVB	5139698	5138957	100.00%	8.749%
4	6.563	757	761	765	rBV	5627807	4877421	94.91%	8.303%
5	6.798	797	801	804	rBV	3052834	2563816	49.89%	4.365%
6	6.951	823	827	831	rBV	4408818	3712891	72.25%	6.321%
7	7.351	889	895	898	rBV	3646718	3085352	60.04%	5.253%
8	7.404	902	904	907	rVV	112135	85661	1.67%	0.146%
9	7.439	907	910	916	rVB4	49137	69433	1.35%	0.118%
10	7.575	928	933	937	rBV4	33846	72940	1.42%	0.124%
11	7.916	987	991	995	rVB6	36352	60740	1.18%	0.103%
12	8.081	1015	1019	1022	rBV	3683732	3336664	64.93%	5.680%
13	8.151	1029	1031	1033	rVB	89650	60809	1.18%	0.104%
14	8.251	1042	1048	1051	rBV7	33945	64451	1.25%	0.110%
15	8.381	1062	1070	1073	rBV7	75960	126594	2.46%	0.216%
16	8.469	1082	1085	1089	rVB3	66837	63115	1.23%	0.107%
17	8.510	1089	1092	1094	rBV	99936	95037	1.85%	0.162%
18	8.586	1102	1105	1108	rVB3	80087	71981	1.40%	0.123%
19	8.628	1108	1112	1117	rBV8	51121	76125	1.48%	0.130%
20	8.681	1118	1121	1124	rBV	207826	169952	3.31%	0.289%
21	8.775	1135	1137	1142	rVB3	65691	95097	1.85%	0.162%
22	9.110	1191	1194	1198	rVB2	107544	94769	1.84%	0.161%
23	9.157	1198	1202	1205	rBV	5244538	4449845	86.59%	7.575%
24	9.245	1214	1217	1221	rBV	256172	235727	4.59%	0.401%
25	9.416	1243	1246	1253	rVB6	40164	58760	1.14%	0.100%
26	9.545	1265	1268	1270	rBV	519737	392058	7.63%	0.667%
27	9.569	1270	1272	1274	rVB	97892	66142	1.29%	0.113%
28	9.775	1305	1307	1311	rVB	231378	160528	3.12%	0.273%
29	9.833	1311	1317	1321	rBV2	3463517	2990423	58.19%	5.091%
30	10.033	1349	1351	1354	rVB	90360	75162	1.46%	0.128%
31	10.098	1359	1362	1366	rBV3	58594	62667	1.22%	0.107%
32	10.275	1389	1392	1394	rVB	194045	140039	2.73%	0.238%
33	10.375	1407	1409	1412	rBV	58456	56718	1.10%	0.097%
34	10.498	1427	1430	1433	rVB	83914	74283	1.45%	0.126%

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35	10.616	1446	1450	1455	rBV	2469598	2040731	39.71%	3.474%
36	10.745	1469	1472	1480	rVB2	209506	282108	5.49%	0.480%
37	11.192	1546	1548	1551	rBV	168527	155446	3.02%	0.265%
38	11.227	1552	1554	1558	rVB	126342	107142	2.08%	0.182%
39	11.316	1565	1569	1571	rBV	2661927	2252548	43.83%	3.835%
40	11.339	1571	1573	1576	rVV	310892	246007	4.79%	0.419%
41	11.386	1579	1581	1584	rVB	84411	77443	1.51%	0.132%
42	11.621	1618	1621	1623	rBV	154082	119310	2.32%	0.203%
43	11.721	1634	1638	1642	rVB	827993	743940	14.48%	1.266%
44	11.851	1656	1660	1665	rBV2	1337520	1169685	22.76%	1.991%
45	11.927	1670	1673	1675	rBV2	130662	137988	2.69%	0.235%
46	12.027	1688	1690	1693	rBV	128756	108004	2.10%	0.184%
47	12.368	1746	1748	1751	rBV2	108485	110246	2.15%	0.188%
48	12.404	1751	1754	1756	rBV	2748070	2351950	45.77%	4.004%
49	12.427	1756	1758	1761	rVB	2790855	2113767	41.13%	3.599%
50	12.527	1770	1775	1778	rBV2	494239	695363	13.53%	1.184%
51	12.557	1778	1780	1790	rVB4	235704	342004	6.66%	0.582%
52	12.633	1790	1793	1799	rBV	720806	835446	16.26%	1.422%
53	12.751	1811	1813	1817	rVB	324930	289513	5.63%	0.493%
54	12.904	1835	1839	1842	rBV	3248413	2866537	55.78%	4.880%
55	13.145	1878	1880	1884	rVB2	181848	195107	3.80%	0.332%
56	13.815	1991	1994	2000	rVB2	630983	801276	15.59%	1.364%
57	13.951	2014	2017	2020	rBV	1781923	1774231	34.53%	3.020%
58	15.398	2259	2263	2266	rVB	1014520	1154462	22.46%	1.965%

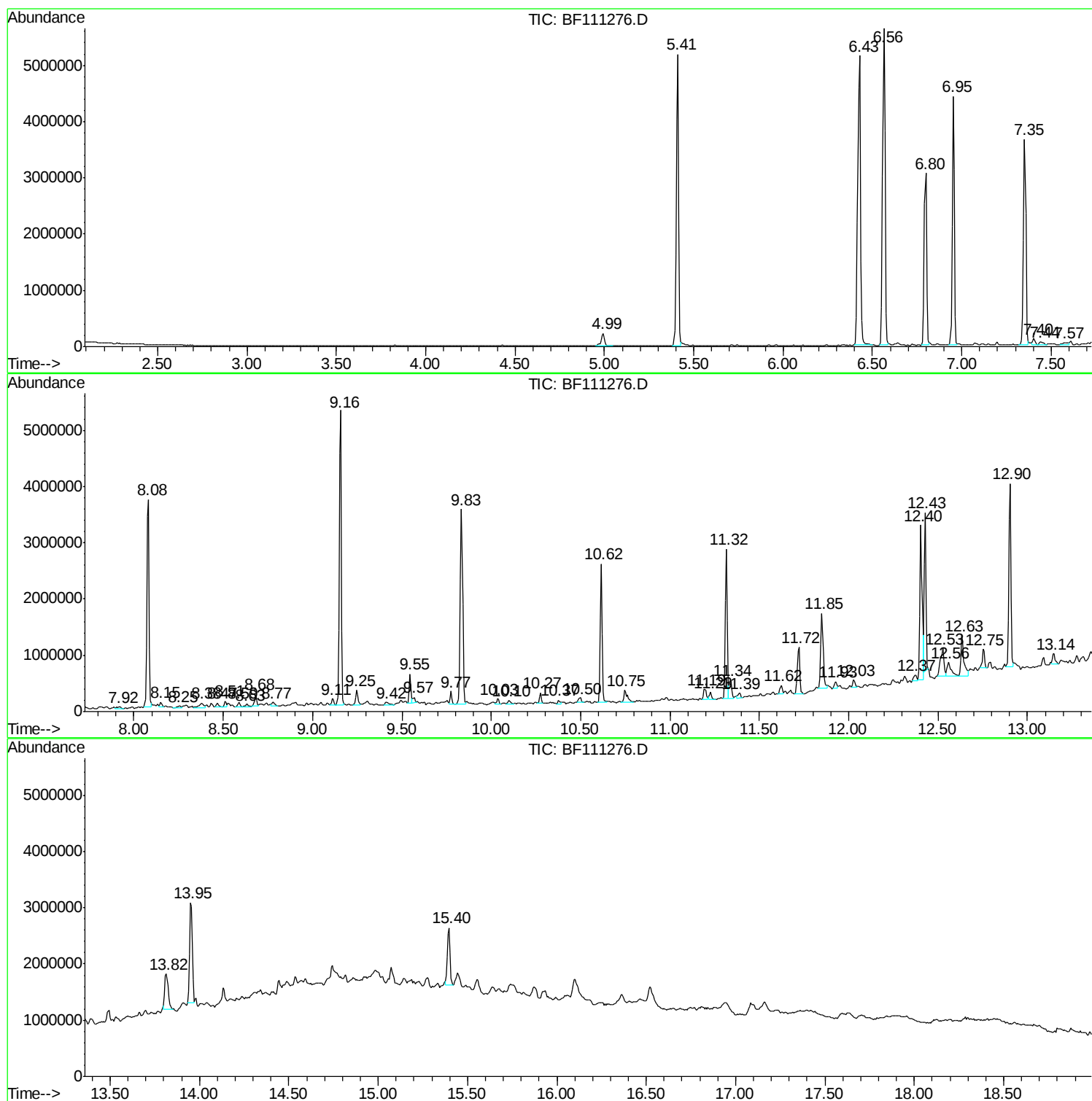
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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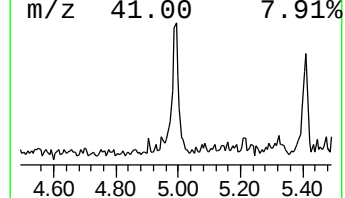
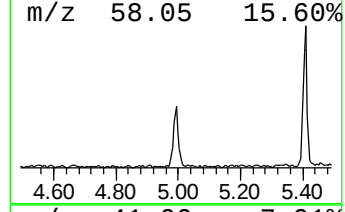
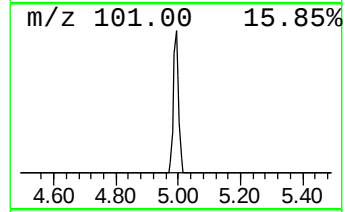
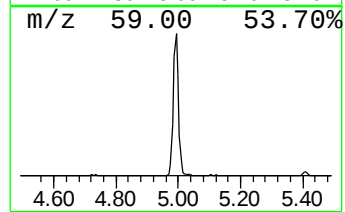
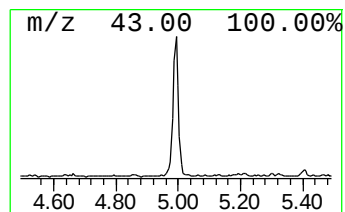
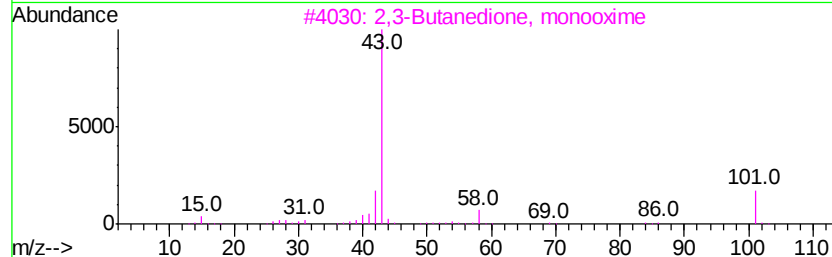
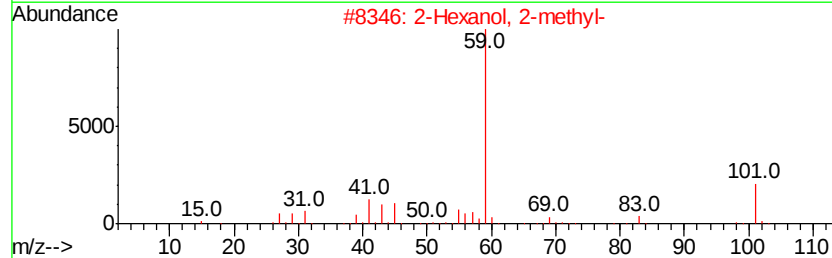
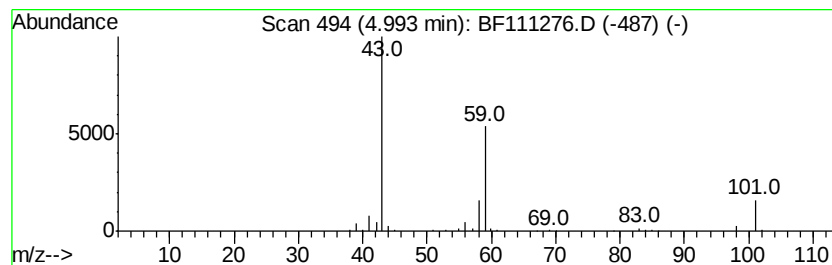
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	2.40 ng	307668	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9



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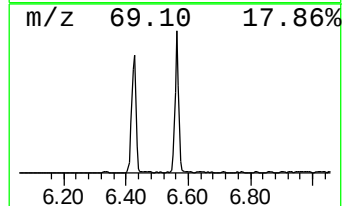
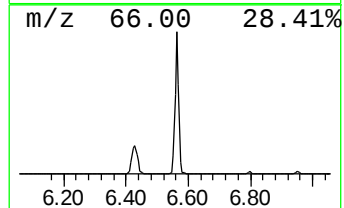
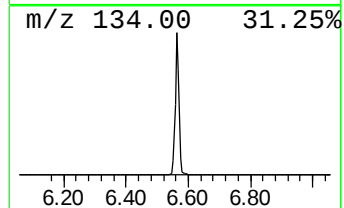
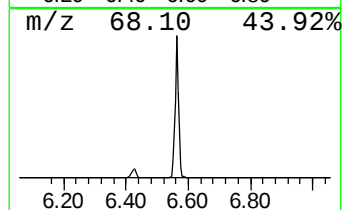
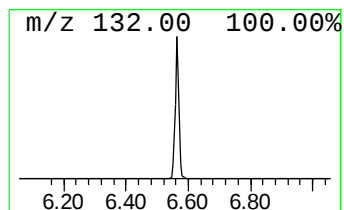
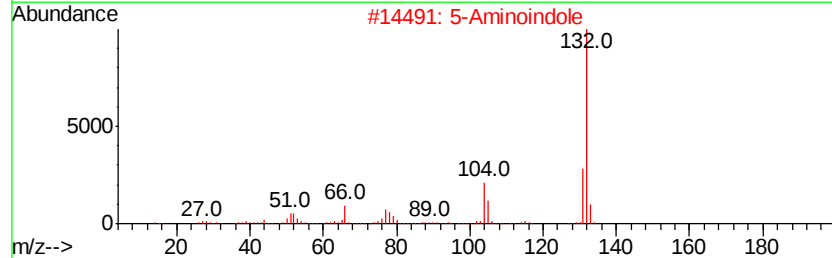
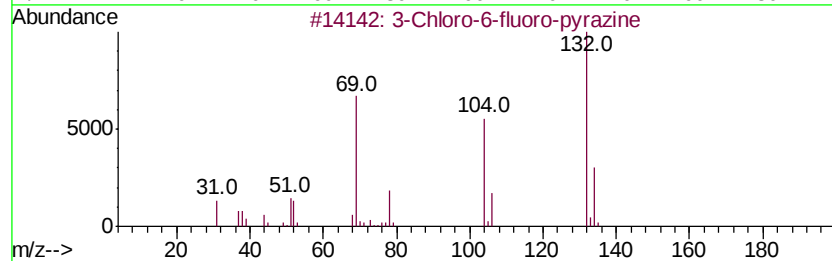
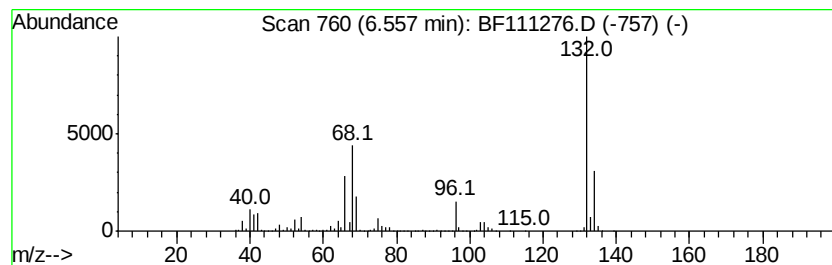
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	38.05 ng	4877420	1,4-Dichlorobenzene-d4	6.80

Hit#	of	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	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9	
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9	



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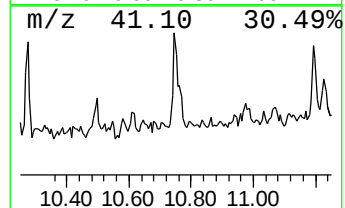
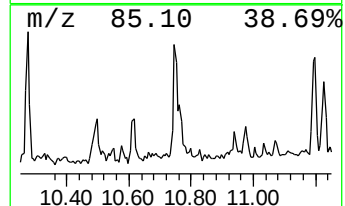
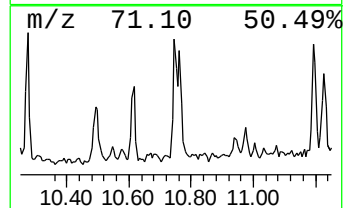
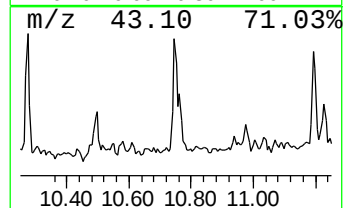
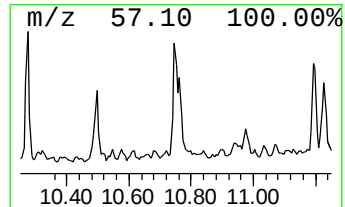
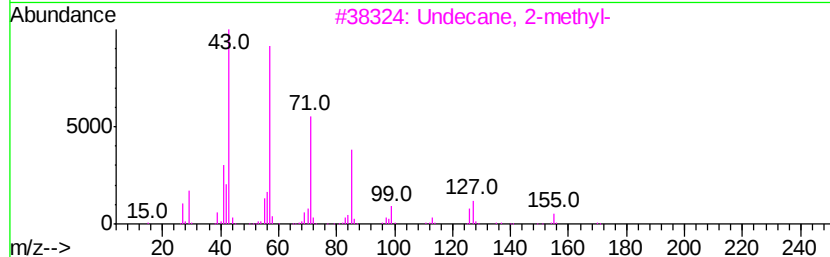
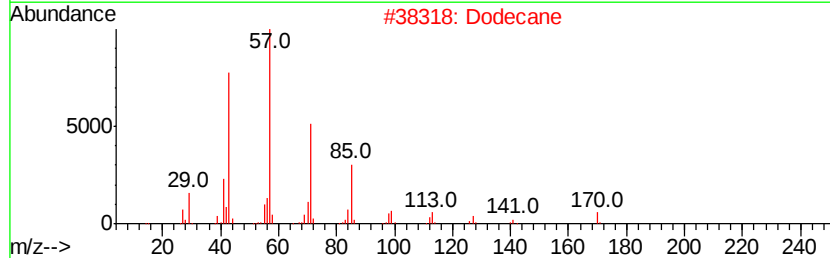
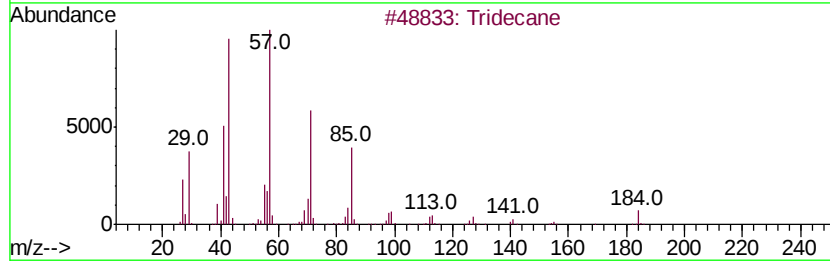
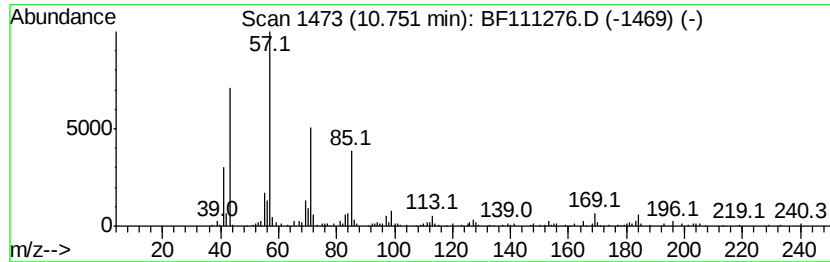
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Tridecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.75	2.50 ng	282108	Phenanthrene-d10		11.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	91
2	Dodecane	170	C12H26	000112-40-3	87
3	Undecane, 2-methyl-	170	C12H26	007045-71-8	80
4	5-Ethyldecane	170	C12H26	017302-36-2	80
5	Tridecane, 2-methyl-	198	C14H30	001560-96-9	74



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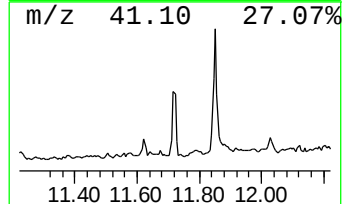
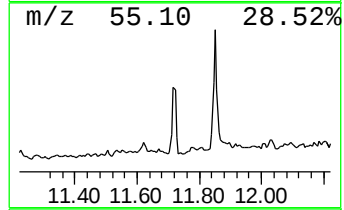
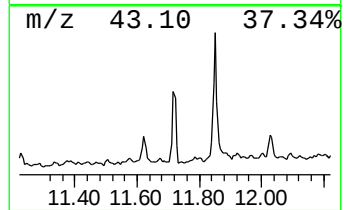
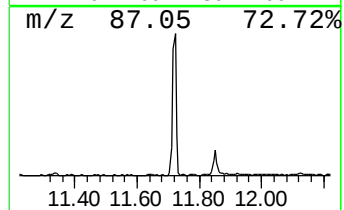
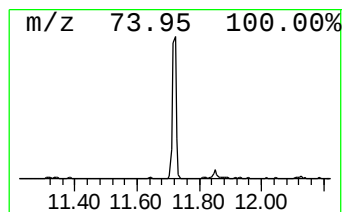
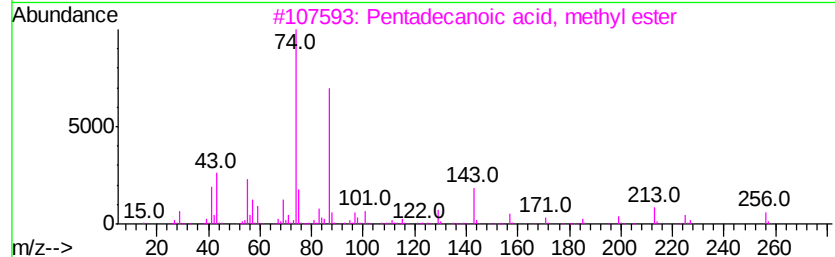
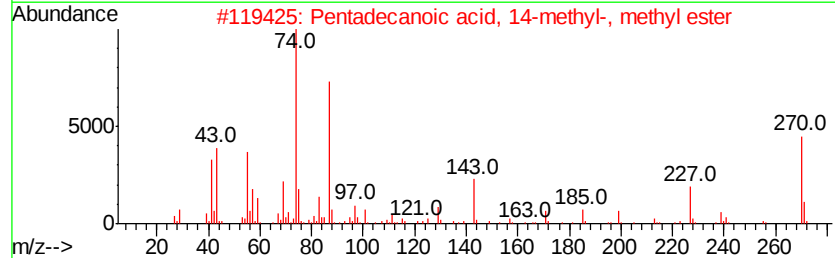
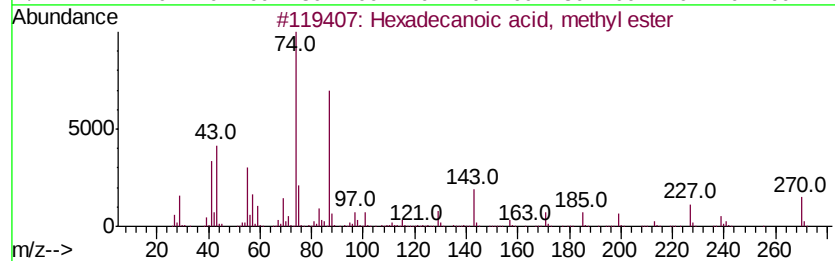
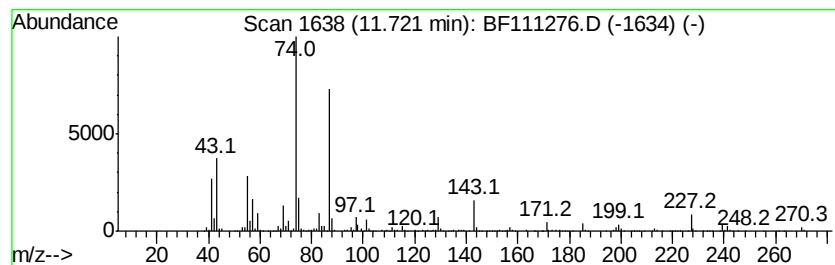
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	6.61 ng	743940	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	86
5		Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	86



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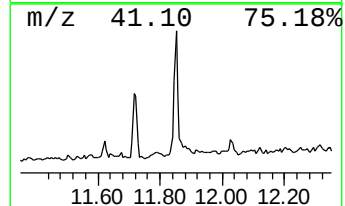
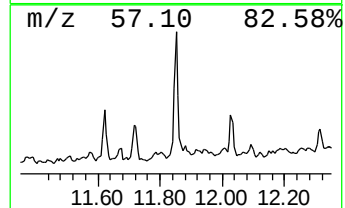
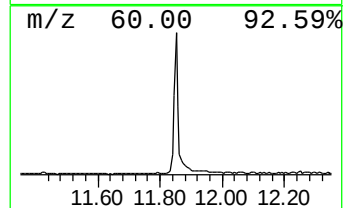
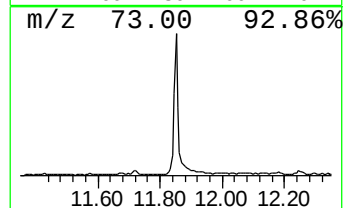
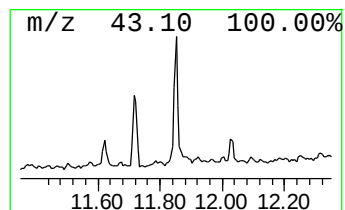
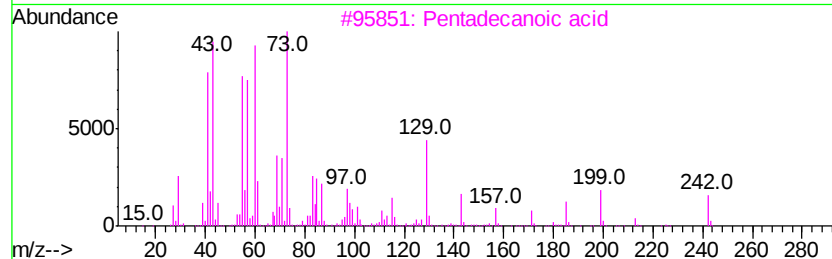
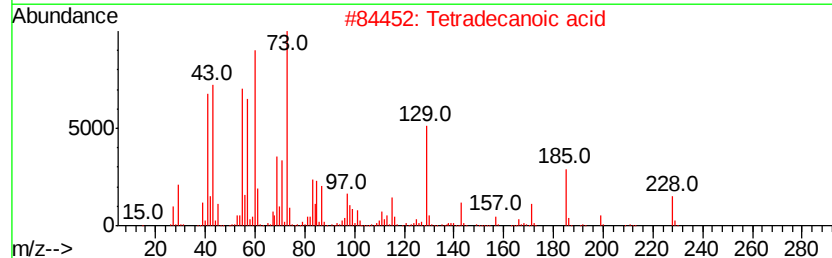
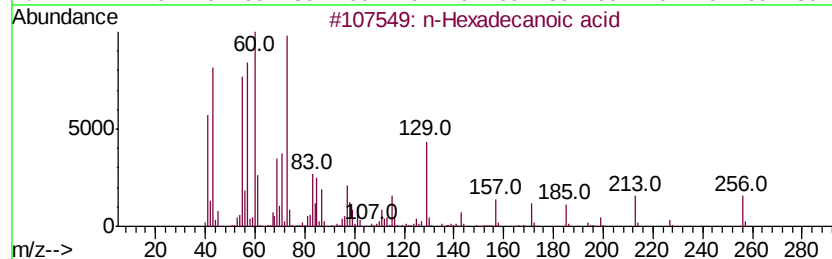
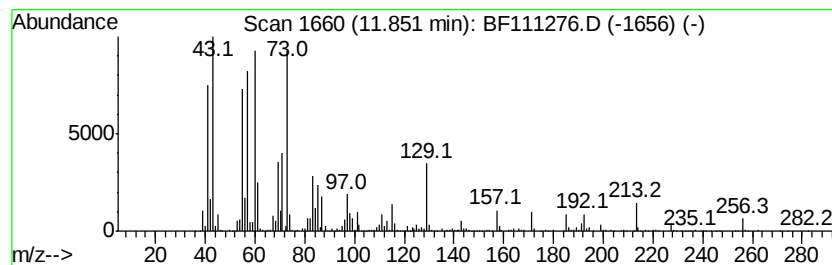
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ClientSampleId :
EO-03-120718-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.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 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.85	10.39 ng	1169690	Phenanthrene-d10		11.32	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid		228	C14H28O2	000544-63-8	93
3	Pentadecanoic acid		242	C15H30O2	001002-84-2	87
4	Undecanoic acid		186	C11H22O2	000112-37-8	72
5	n-Decanoic acid		172	C10H20O2	000334-48-5	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
Data File : BF111276.D
Acq On : 11 Dec 2018 14:04
Operator : JU/SJ
Sample : J6196-01 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-03-120718-A

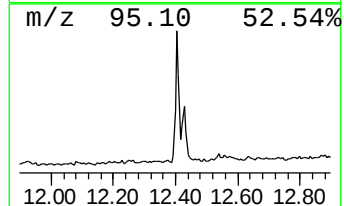
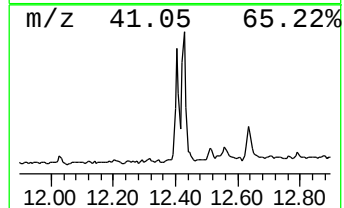
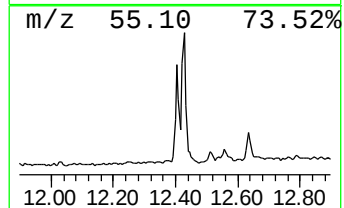
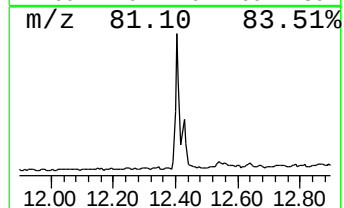
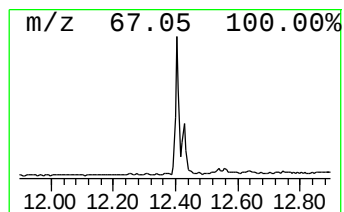
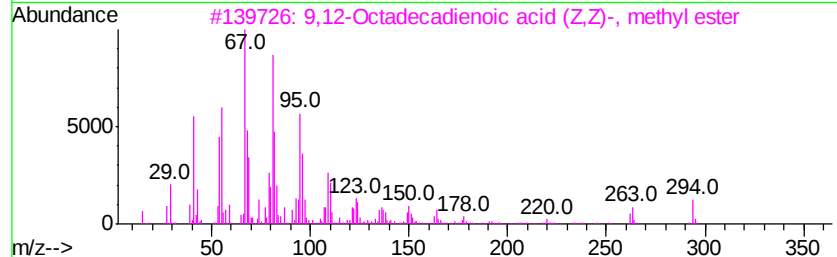
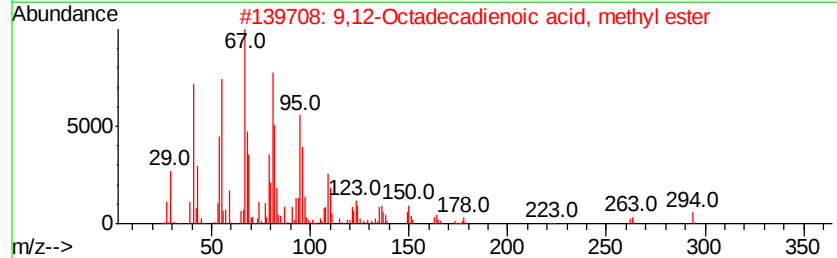
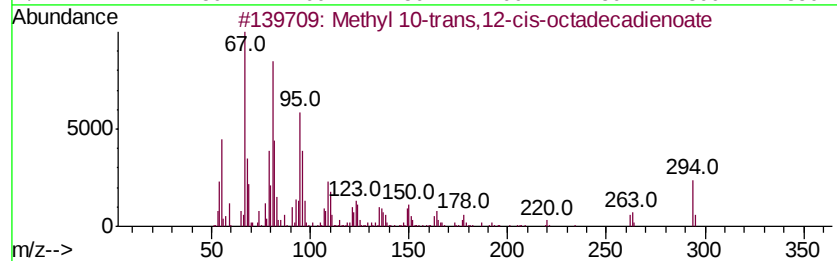
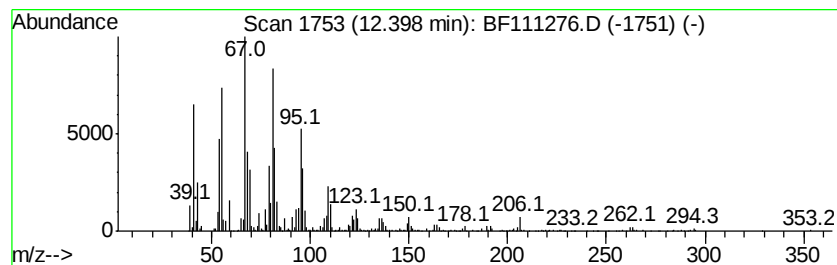
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Methyl 10-trans,12-cis-octa... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	20.88 ng	2351950	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99
2		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
3		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
5		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
Data File : BF111276.D
Acq On : 11 Dec 2018 14:04
Operator : JU/SJ
Sample : J6196-01 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-03-120718-A

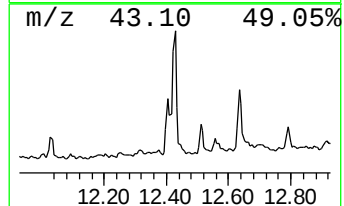
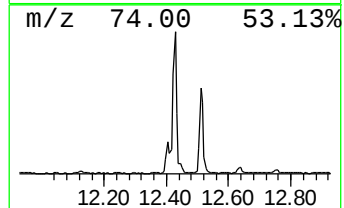
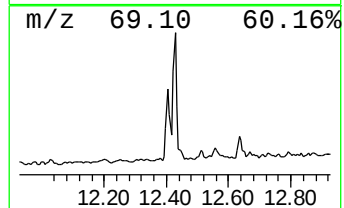
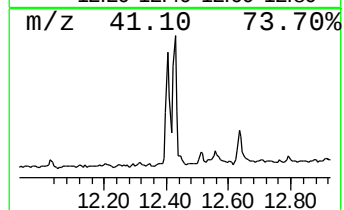
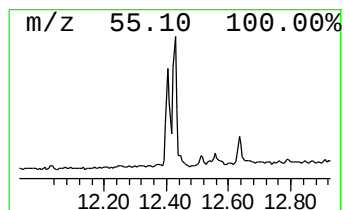
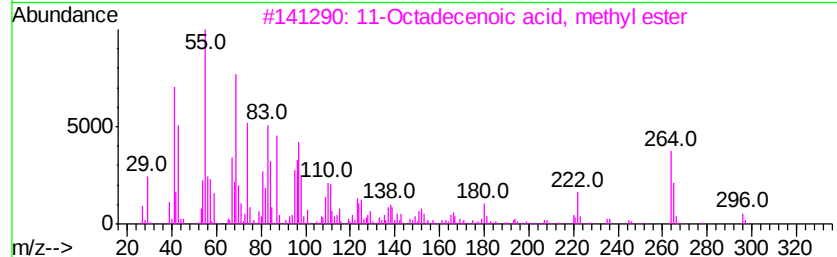
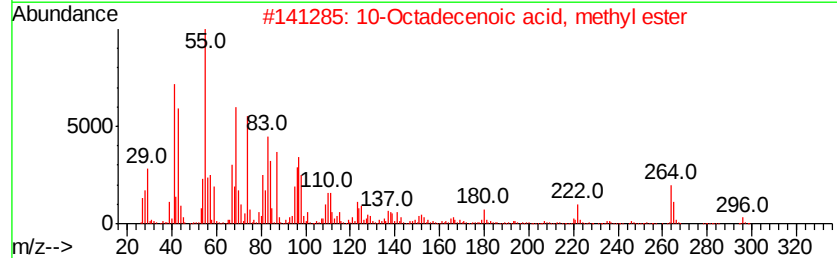
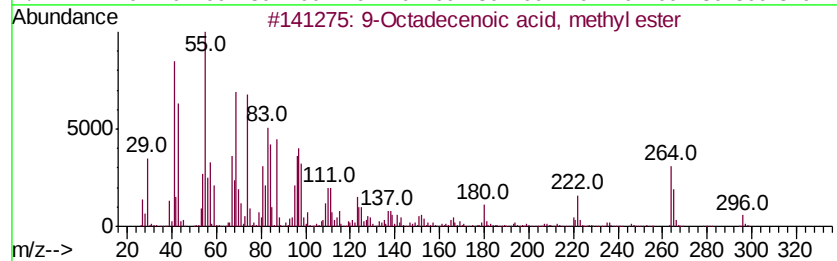
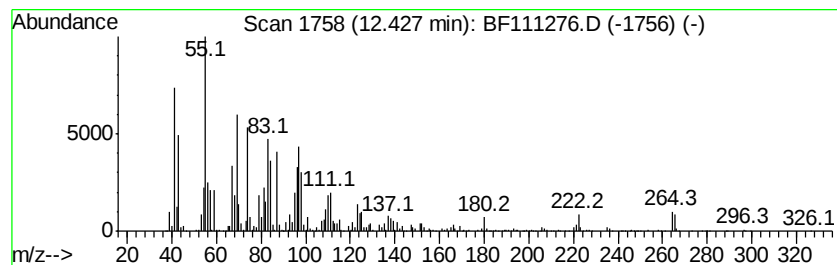
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9-Octadecenoic acid, methyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	18.77 ng	2113770	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
2		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
3		11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	99
4		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
5		cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1000333-58-3	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
Data File : BF111276.D
Acq On : 11 Dec 2018 14:04
Operator : JU/SJ
Sample : J6196-01 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-03-120718-A

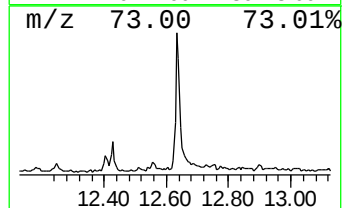
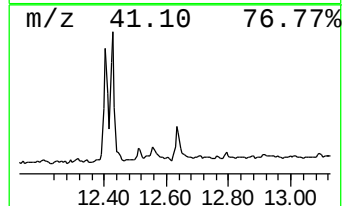
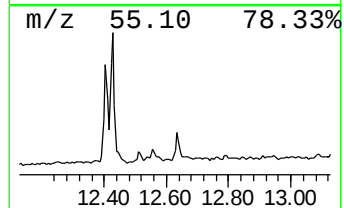
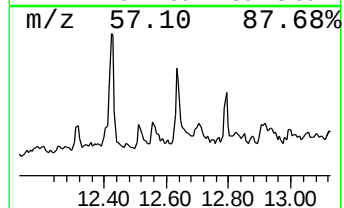
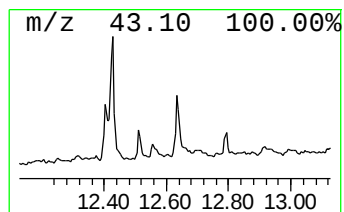
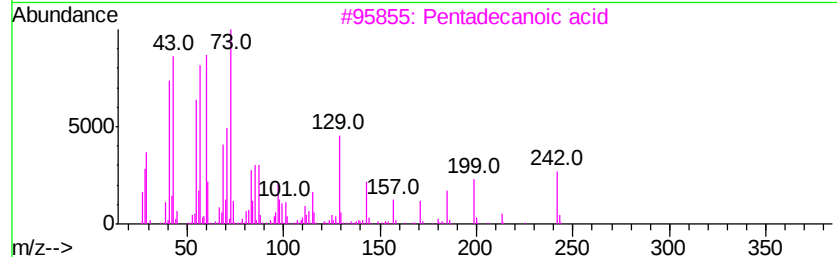
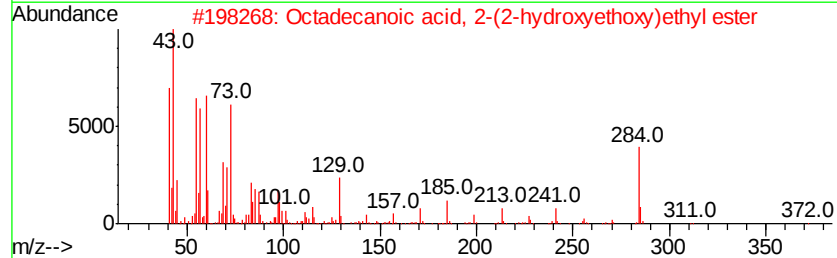
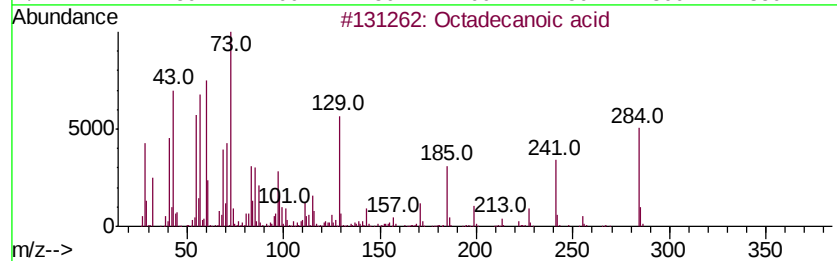
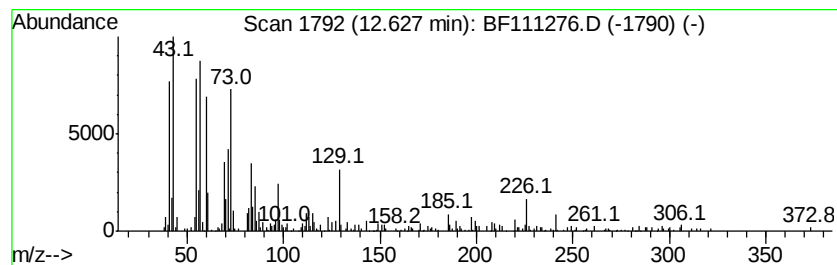
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	7.42 ng	835446	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	95
2		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	62
4		Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	58
5		Undecanoic acid	186	C11H22O2	000112-37-8	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
Data File : BF111276.D
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Sample : J6196-01 2X
Misc :
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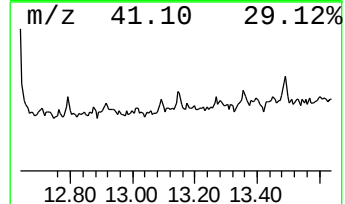
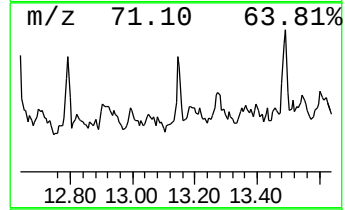
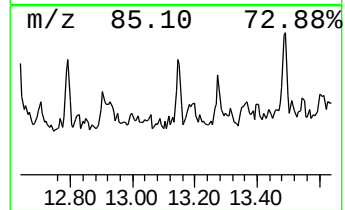
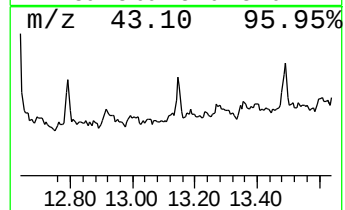
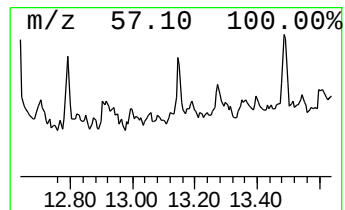
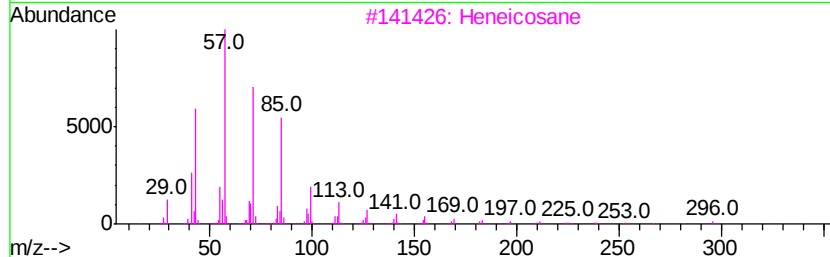
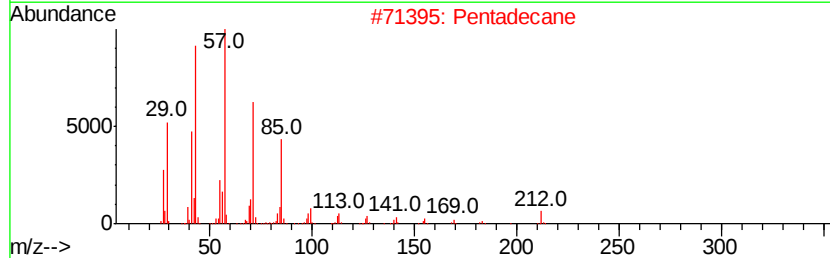
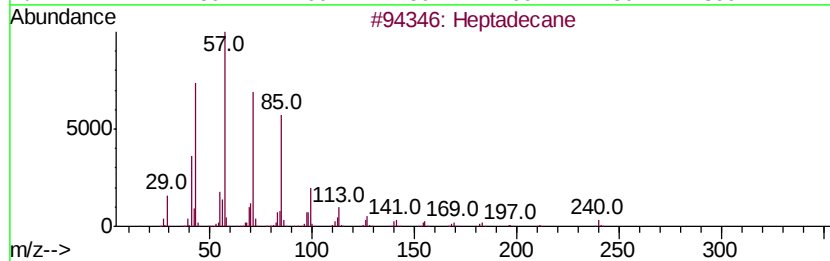
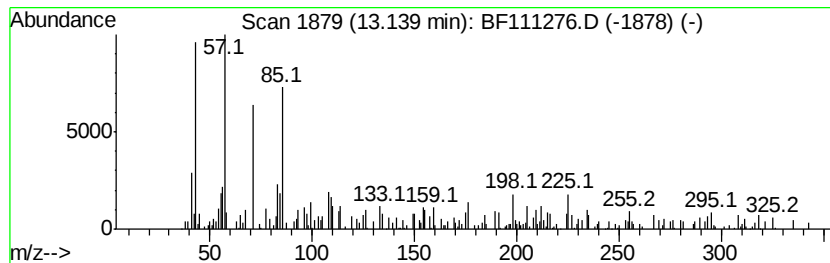
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.14	2.20 ng	195107	Chrysene-d12	13.96	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	55
2	Pentadecane	212	C15H32	000629-62-9	55
3	Heneicosane	296	C21H44	000629-94-7	45
4	Eicosane	282	C20H42	000112-95-8	45
5	2-Bromo dodecane	248	C12H25Br	013187-99-0	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
Data File : BF111276.D
Acq On : 11 Dec 2018 14:04
Operator : JU/SJ
Sample : J6196-01 2X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-03-120718-A

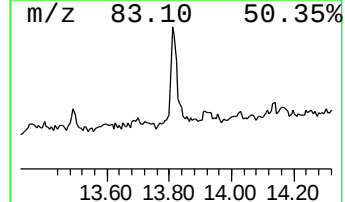
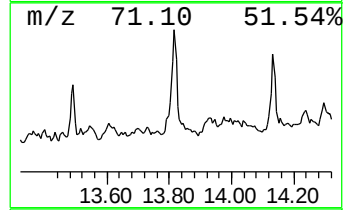
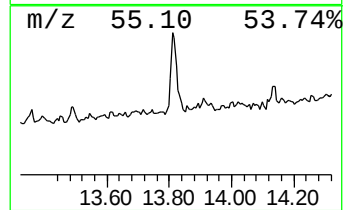
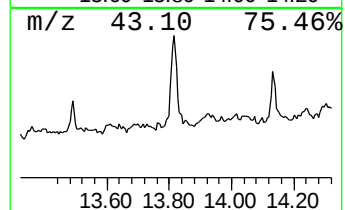
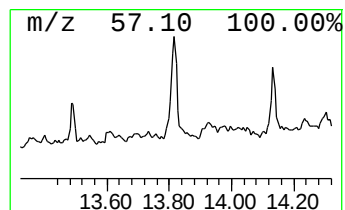
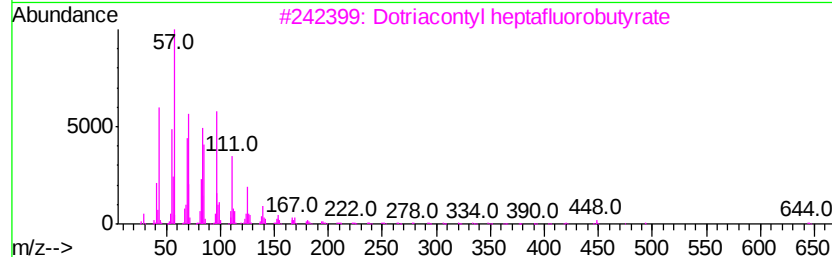
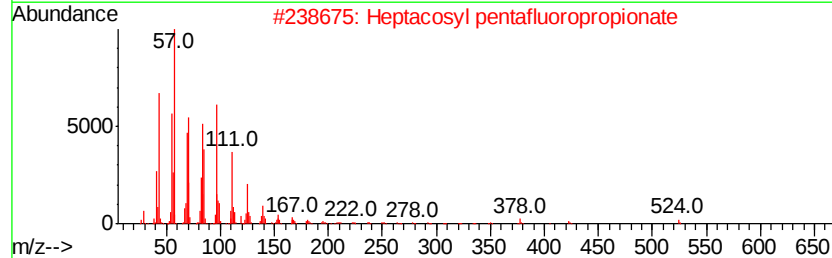
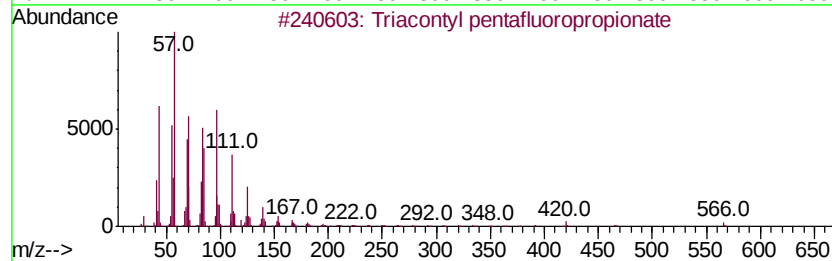
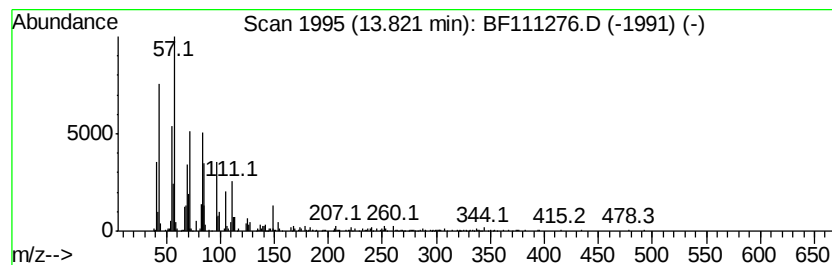
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triacontyl pentafluoropropi... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	9.03 ng	801276	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontyl pentafluoropropionate	584	C33H61F5O2	1000351-80-0	91
2		Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	91
3		Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	91
4		Dotriacontyl trifluoroacetate	562	C34H65F3O2	1000351-75-4	91
5		Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121118\
 Data File : BF111276.D
 Acq On : 11 Dec 2018 14:04
 Operator : JU/SJ
 Sample : J6196-01 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120818.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...	4.99	2.4	ng	307668	1	6.80	2563820	20.0
unknown6.56	6.56	38.0	ng	4877420	1	6.80	2563820	20.0
Tridecane	10.75	2.5	ng	282108	4	11.32	2252550	20.0
Hexadecanoic acid...	11.72	6.6	ng	743940	4	11.32	2252550	20.0
n-Hexadecanoic acid	11.85	10.4	ng	1169690	4	11.32	2252550	20.0
Methyl 10-trans,1...	12.40	20.9	ng	2351950	4	11.32	2252550	20.0
9-Octadecenoic ac...	12.43	18.8	ng	2113770	4	11.32	2252550	20.0
Octadecanoic acid	12.63	7.4	ng	835446	4	11.32	2252550	20.0
Heptadecane	13.14	2.2	ng	195107	5	13.96	1774230	20.0
Triacontyl pentaf...	13.82	9.0	ng	801276	5	13.96	1774230	20.0