

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109401.D
 Acq On : 27 Sep 2018 21:28
 Operator : JU/SJ
 Sample : J5077-06
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PARCEL-193-092418-2

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-BF092218.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	2.440	55	60	66	rBV	71027	104258	4.97%	0.484%
2	4.899	475	478	487	rBV	59063	75256	3.59%	0.349%
3	5.316	545	549	553	rBV	1778097	1813757	86.51%	8.414%
4	6.345	719	724	727	rBV	1911777	1932611	92.18%	8.966%
5	6.475	742	746	750	rBV	2149978	1946202	92.83%	9.029%
6	6.704	782	785	789	rBV	739316	641469	30.60%	2.976%
7	6.863	808	812	815	rBV	1894812	1531394	73.04%	7.105%
8	7.269	876	881	884	rBV	1319157	1213218	57.87%	5.628%
9	7.934	991	994	999	rBV	36428	33775	1.61%	0.157%
10	7.987	999	1003	1009	rBV	992975	792596	37.80%	3.677%
11	9.063	1182	1186	1189	rBV	2529697	2096628	100.00%	9.727%
12	9.457	1249	1253	1256	rBV	765517	578563	27.59%	2.684%
13	9.598	1274	1277	1283	rVB	35601	29608	1.41%	0.137%
14	9.739	1297	1301	1304	rBV	848407	785065	37.44%	3.642%
15	10.522	1430	1434	1439	rBV	1126792	1030373	49.14%	4.780%
16	11.216	1548	1552	1554	rBV	711080	675827	32.23%	3.135%
17	11.233	1554	1555	1558	rVV	155977	116550	5.56%	0.541%
18	11.286	1561	1564	1568	rVB	45853	37567	1.79%	0.174%
19	11.457	1586	1593	1601	rBV3	35073	80836	3.86%	0.375%
20	11.716	1634	1637	1639	rBV	52080	45436	2.17%	0.211%
21	11.751	1639	1643	1647	rVV2	122958	140251	6.69%	0.651%
22	11.822	1652	1655	1658	rVV	70877	76781	3.66%	0.356%
23	11.845	1658	1659	1664	rVB	41013	28571	1.36%	0.133%
24	12.027	1688	1690	1697	rVB2	30244	32027	1.53%	0.149%
25	12.263	1726	1730	1733	rBV	53787	62122	2.96%	0.288%
26	12.345	1741	1744	1748	rVB2	27925	33223	1.58%	0.154%
27	12.422	1753	1757	1760	rBV	262895	246426	11.75%	1.143%
28	12.645	1789	1795	1798	rBV	315324	290289	13.85%	1.347%
29	12.704	1803	1805	1809	rVB3	24163	25733	1.23%	0.119%
30	12.792	1817	1820	1824	rVV	1673039	1582241	75.47%	7.340%
31	12.827	1824	1826	1829	rVB	85738	68313	3.26%	0.317%
32	12.869	1829	1833	1836	rBV2	37075	47366	2.26%	0.220%
33	12.951	1844	1847	1849	rBV4	23951	30354	1.45%	0.141%
34	12.974	1849	1851	1855	rVB	57707	53141	2.53%	0.247%

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35	13.039	1857	1862	1865	rBV	55743	61778	2.95%	0.287%
36	13.080	1865	1869	1872	rVV2	49725	50509	2.41%	0.234%
37	13.169	1881	1884	1887	rVB	41518	32393	1.55%	0.150%
38	13.198	1887	1889	1893	rVB	41714	34043	1.62%	0.158%
39	13.380	1917	1920	1925	rVB3	30011	38211	1.82%	0.177%
40	13.480	1934	1937	1941	rVB2	29913	42109	2.01%	0.195%
41	13.586	1951	1955	1958	rBV3	41426	62139	2.96%	0.288%
42	13.710	1973	1976	1982	rBV2	96141	126543	6.04%	0.587%
43	13.839	1993	1998	2001	rBV	898516	907958	43.31%	4.212%
44	13.863	2001	2002	2005	rVB	156125	90749	4.33%	0.421%
45	14.021	2027	2029	2032	rVB3	34420	30340	1.45%	0.141%
46	14.257	2067	2069	2073	rBV	29689	27196	1.30%	0.126%
47	14.321	2077	2080	2082	rBV3	64842	64807	3.09%	0.301%
48	14.345	2082	2084	2086	rBV2	36870	34826	1.66%	0.162%
49	14.621	2128	2131	2133	rBV	43171	39811	1.90%	0.185%
50	14.686	2139	2142	2147	rVB	184106	181320	8.65%	0.841%
51	14.751	2150	2153	2158	rVB	164252	166414	7.94%	0.772%
52	14.845	2166	2169	2176	rVB	116823	198036	9.45%	0.919%
53	14.933	2181	2184	2187	rVV2	101796	105841	5.05%	0.491%
54	14.968	2187	2190	2198	rVB5	58937	116461	5.55%	0.540%
55	15.121	2214	2216	2222	rVB	85481	97945	4.67%	0.454%
56	15.180	2223	2226	2230	rVB	79106	89306	4.26%	0.414%
57	15.239	2232	2236	2240	rBV	392596	448278	21.38%	2.080%
58	15.468	2272	2275	2278	rVB	65871	73552	3.51%	0.341%
59	15.686	2308	2312	2317	rBV2	111918	156777	7.48%	0.727%

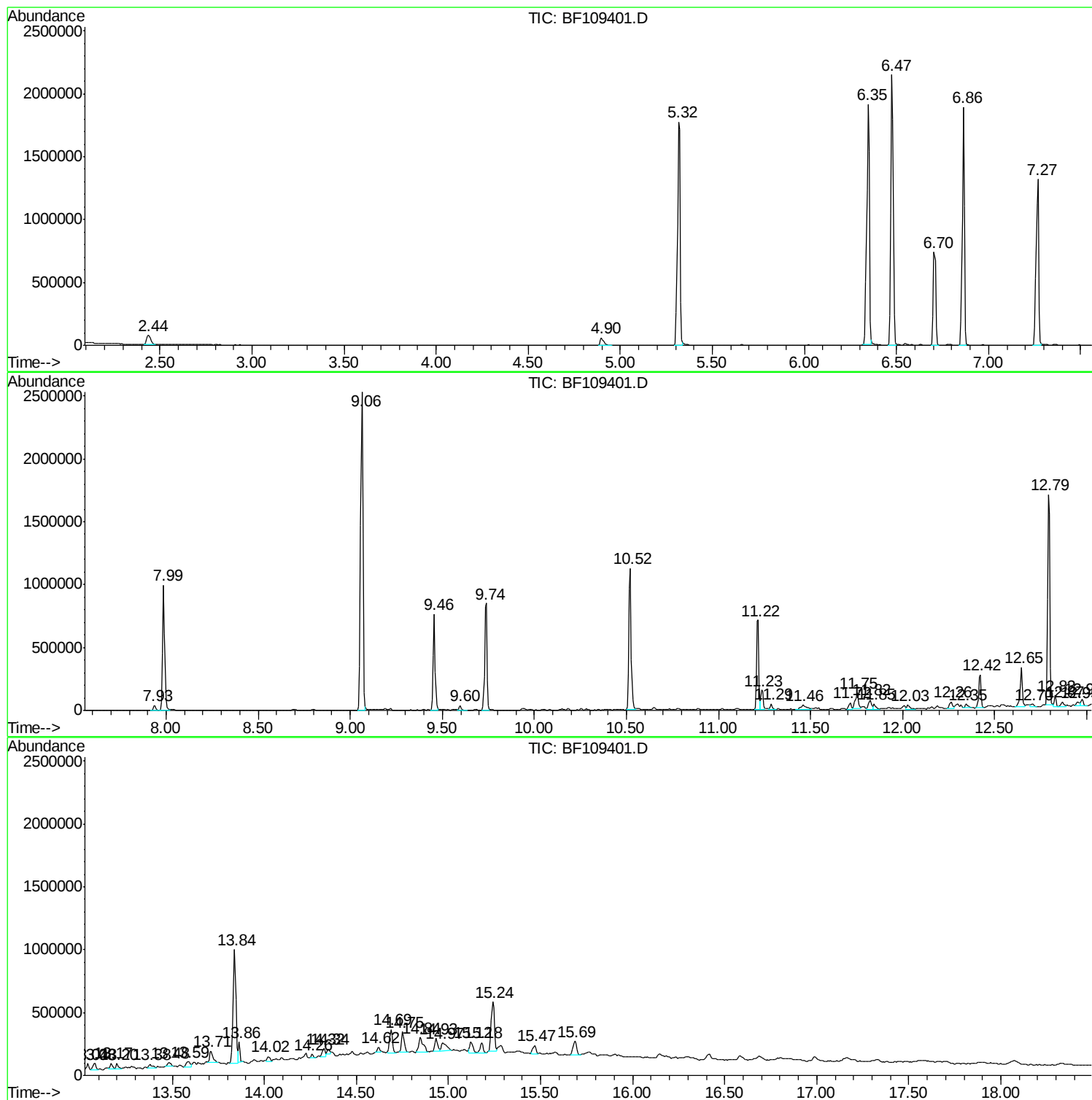
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Operator : JU/SJ
Sample : J5077-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

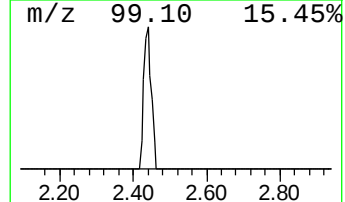
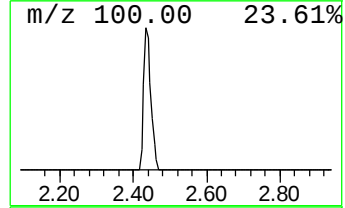
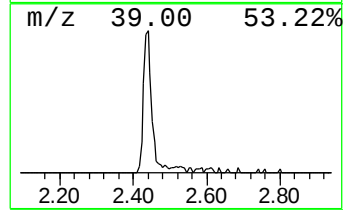
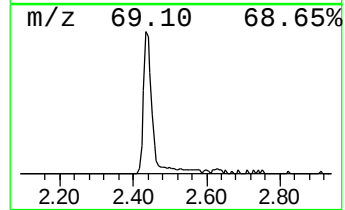
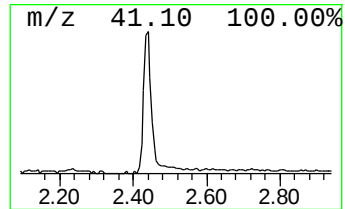
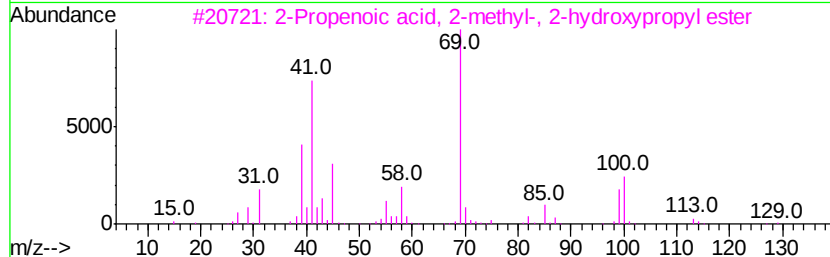
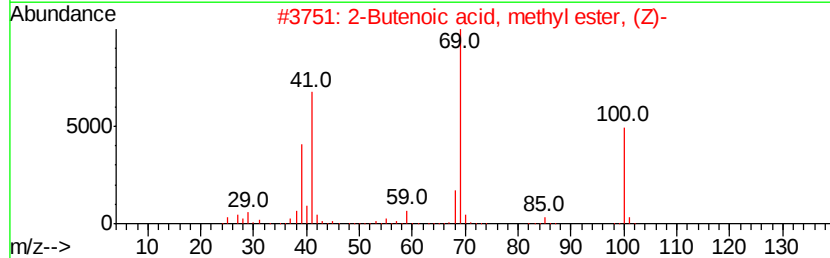
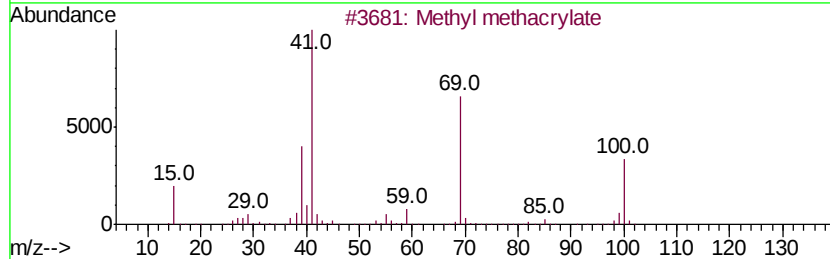
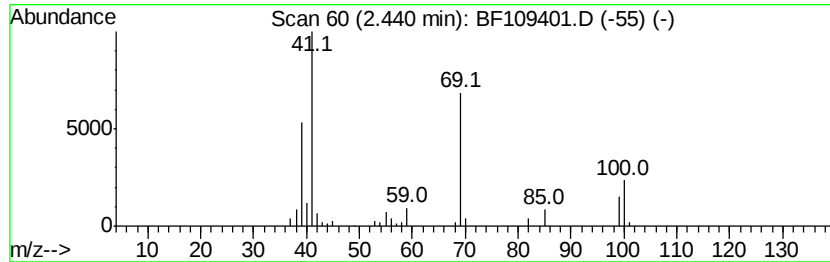
Instrument :
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ClientSampleId :
PARCEL-193-092418-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Methyl methacrylate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
2.44	3.25 ng	104258	1,4-Dichlorobenzene-d4		6.70	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl methacrylate	100	C5H8O2	000080-62-6	80
2		2-Butenoic acid, methyl ester, (Z)-	100	C5H8O2	004358-59-2	80
3		2-Propenoic acid, 2-methyl-, 2-h...	144	C7H12O3	000923-26-2	56
4		2-Butenoic acid, methyl ester, (E)-	100	C5H8O2	000623-43-8	53
5		Methyl 3-butenate	100	C5H8O2	003724-55-8	50



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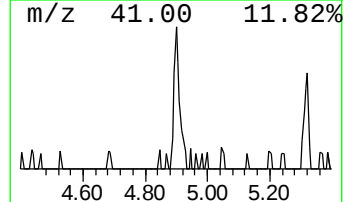
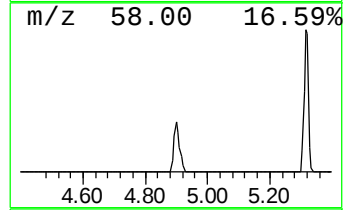
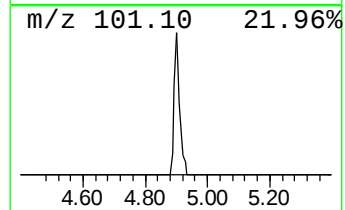
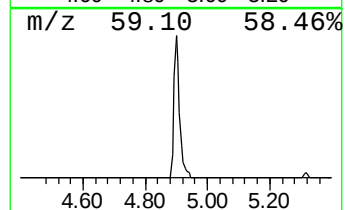
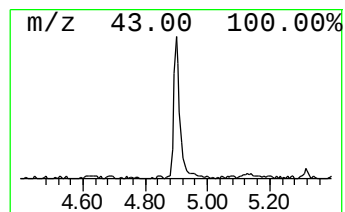
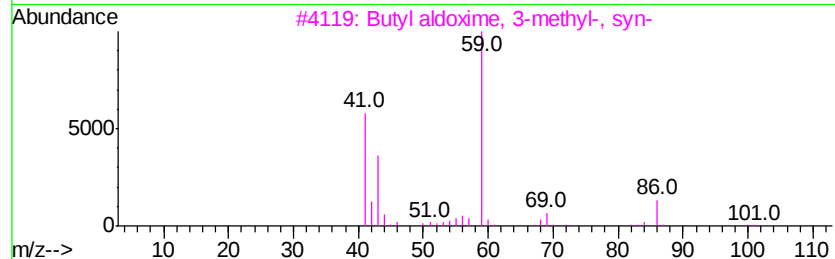
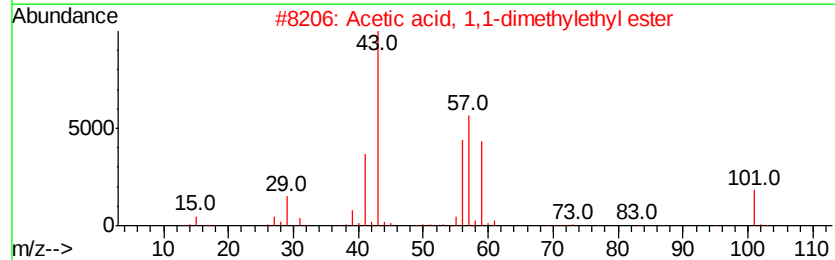
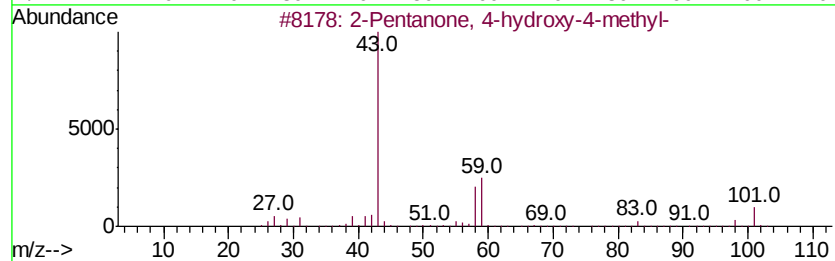
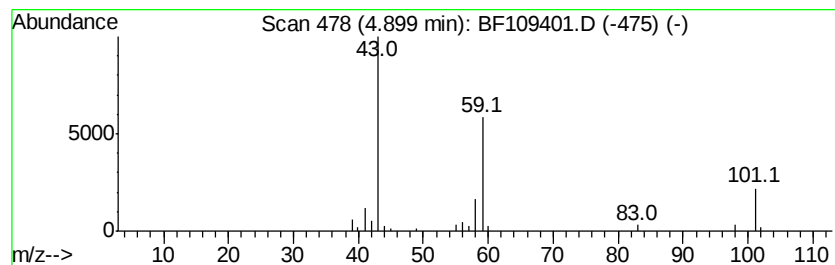
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	2.35 ng	75256	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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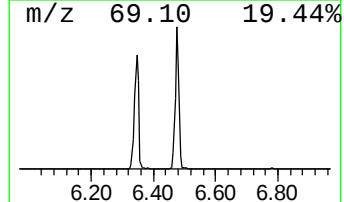
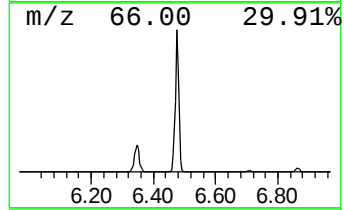
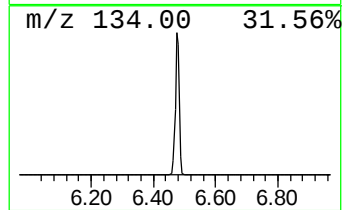
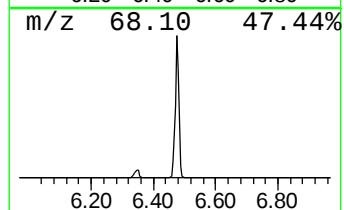
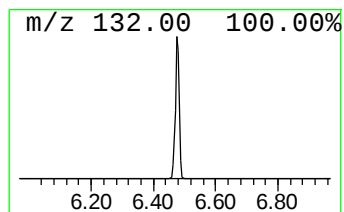
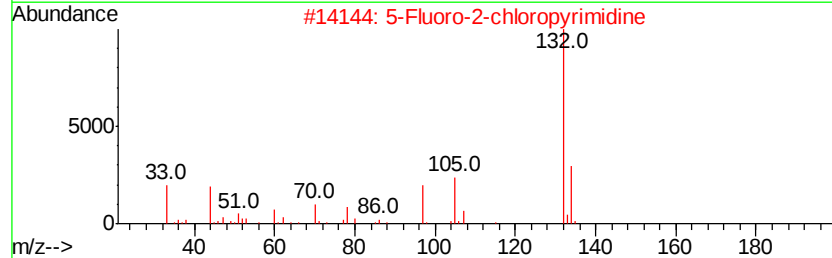
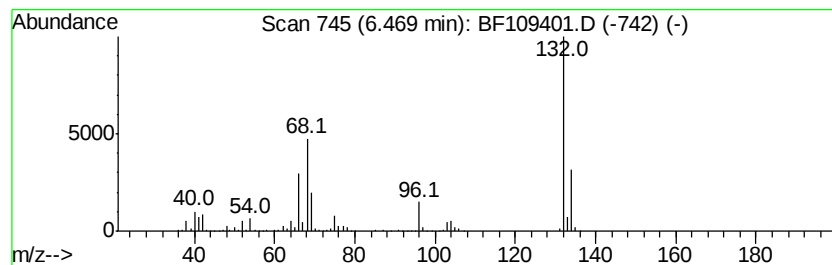
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	60.68 ng	1946200	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16	
3	Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10	
4	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9	
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9	



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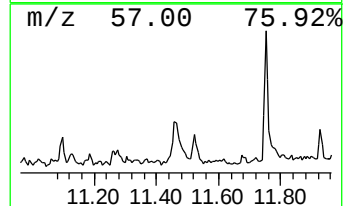
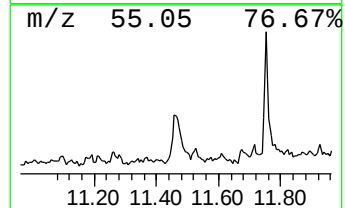
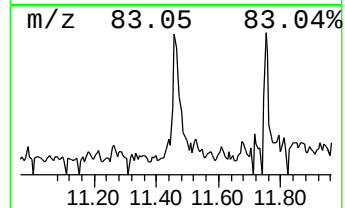
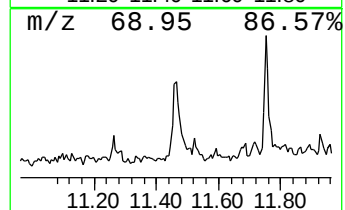
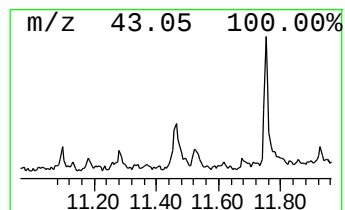
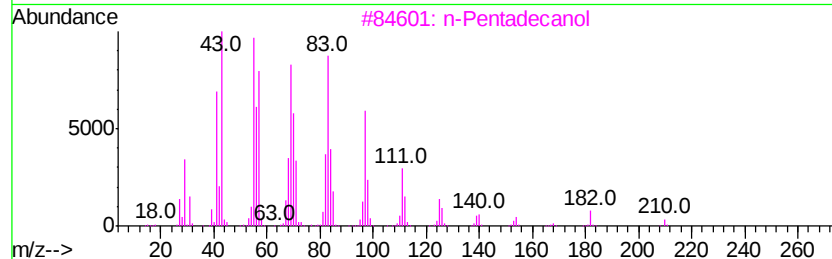
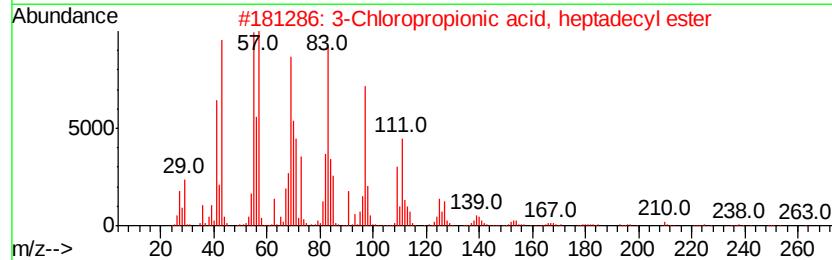
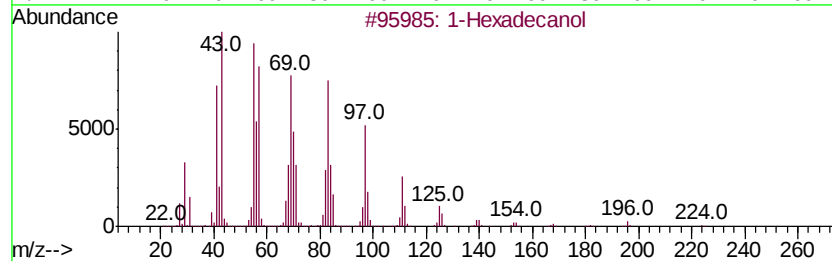
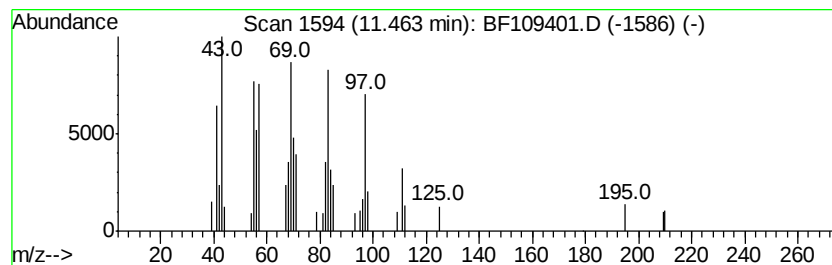
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Hexadecanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.46	2.39 ng	80836	Phenanthrene-d10	11.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexadecanol	242	C16H34O	036653-82-4	91
2		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91
3		n-Pentadecanol	228	C15H32O	000629-76-5	91
4		n-Heptadecanol-1	256	C17H36O	001454-85-9	91
5		n-Tridecan-1-ol	200	C13H28O	000112-70-9	90



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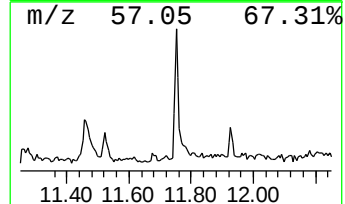
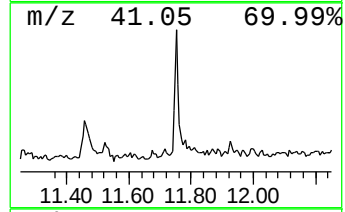
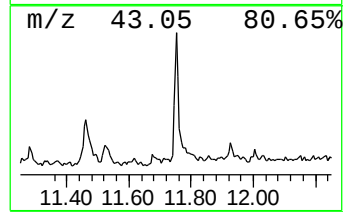
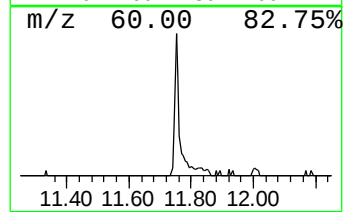
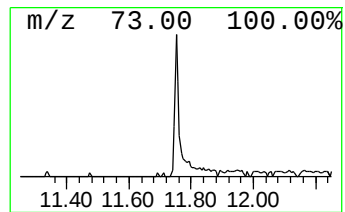
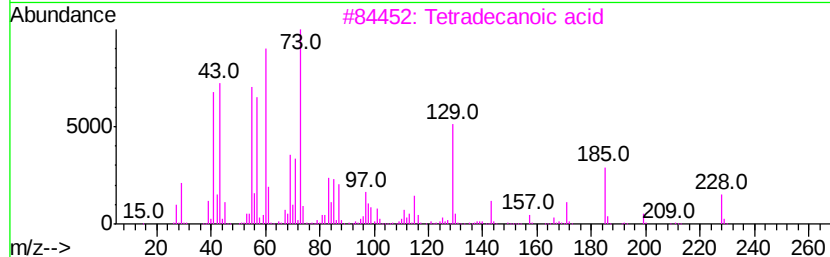
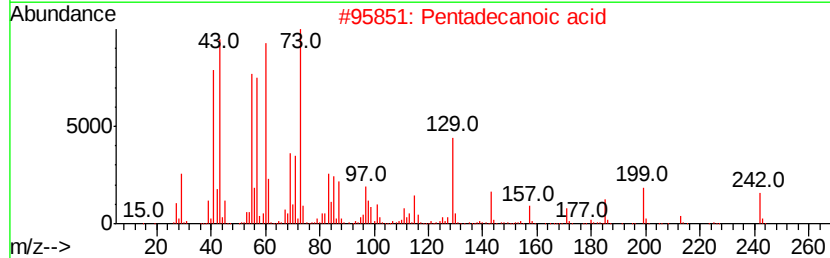
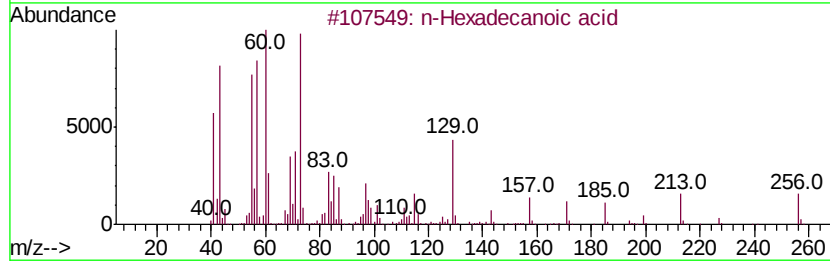
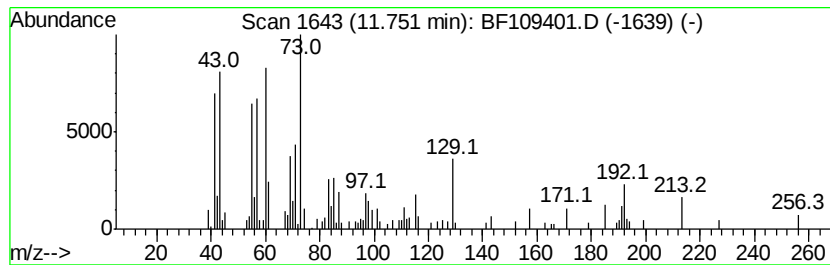
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	4.15 ng	140251	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	70
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	58
4		Tridecanoic acid	214	C13H26O2	000638-53-9	58
5		1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
Data File : BF109401.D
Acq On : 27 Sep 2018 21:28
Operator : JU/SJ
Sample : J5077-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PARCEL-193-092418-2

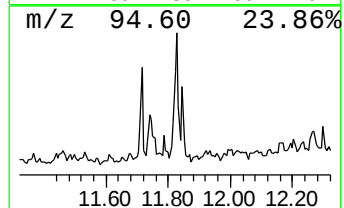
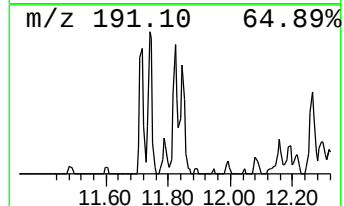
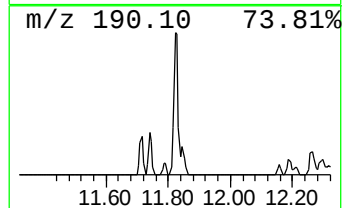
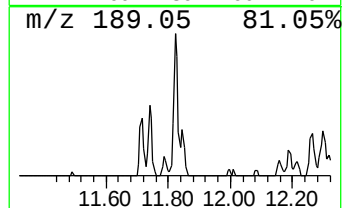
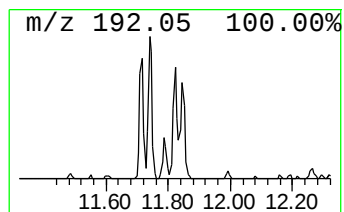
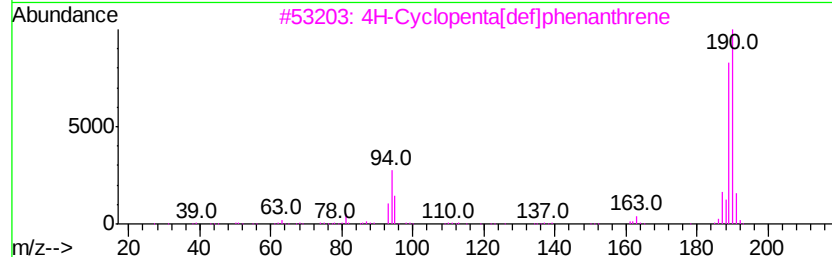
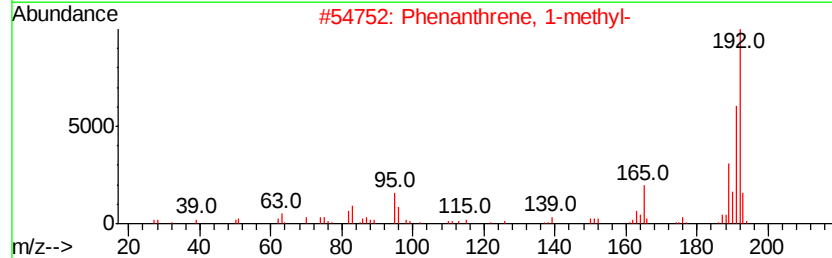
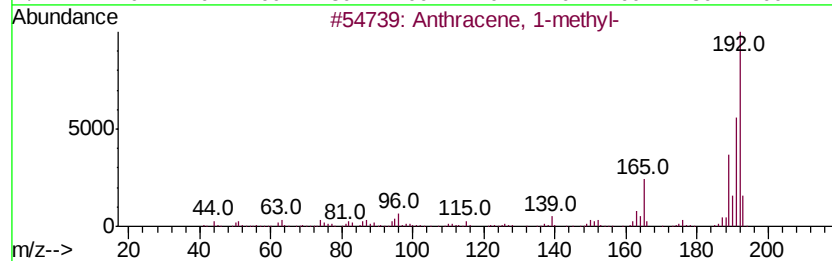
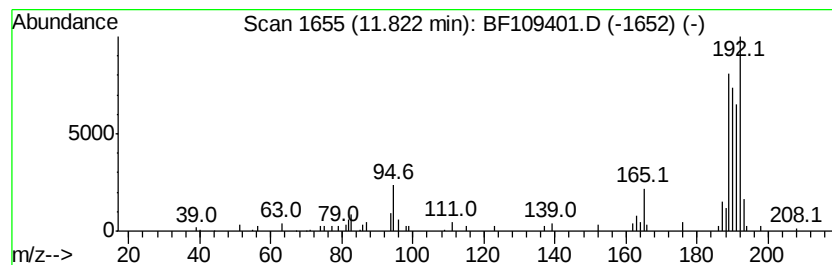
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Anthracene, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	2.27 ng	76781	Phenanthrene-d10	11.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	80
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	60
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
4		8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	51
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
Data File : BF109401.D
Acq On : 27 Sep 2018 21:28
Operator : JU/SJ
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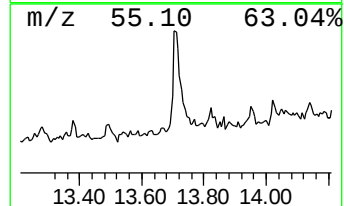
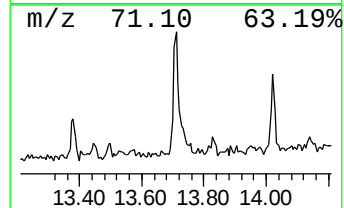
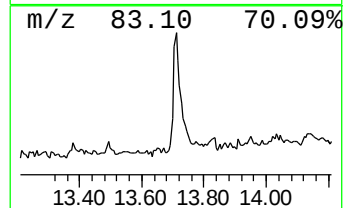
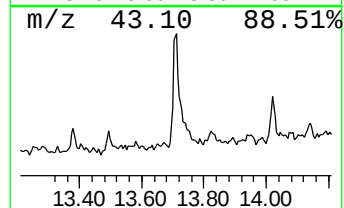
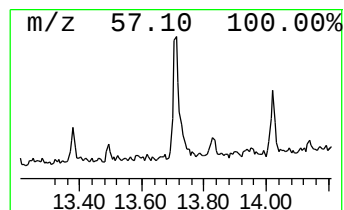
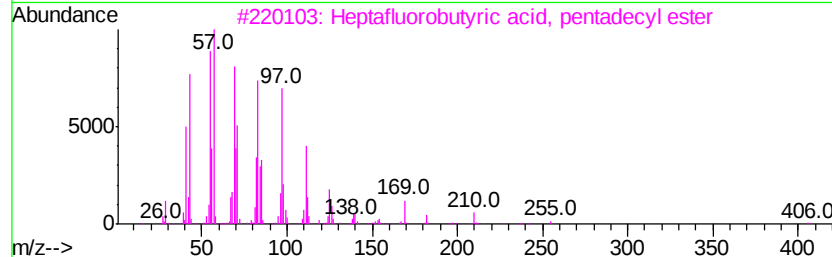
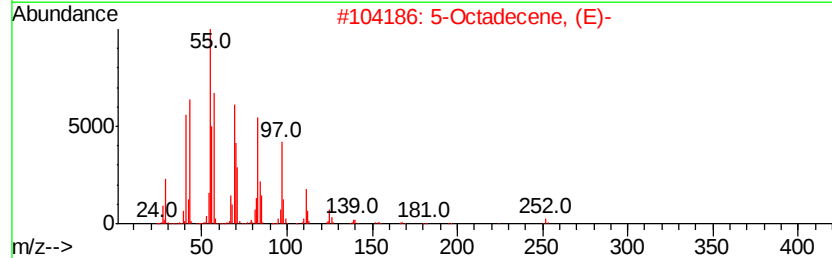
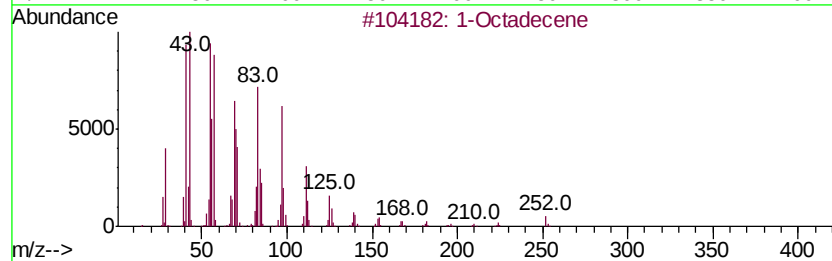
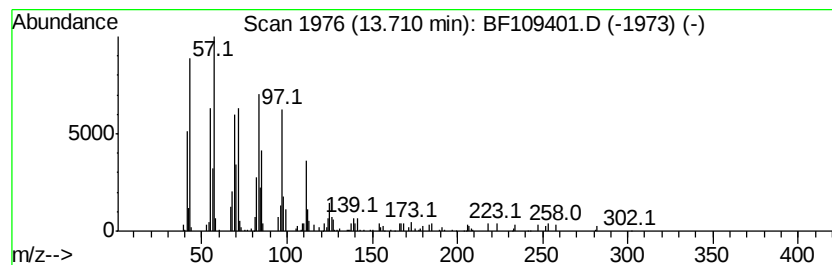
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Octadecene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	2.79 ng	126543	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecene	252	C18H36	000112-88-9	93
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	93
3		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
4		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
Data File : BF109401.D
Acq On : 27 Sep 2018 21:28
Operator : JU/SJ
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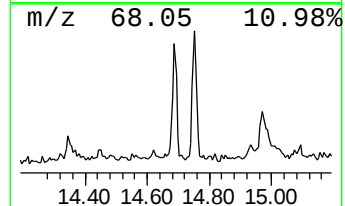
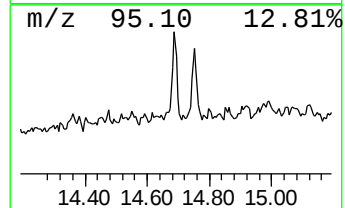
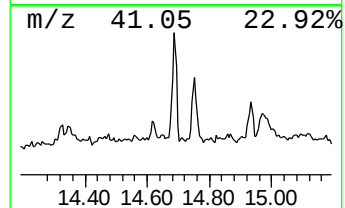
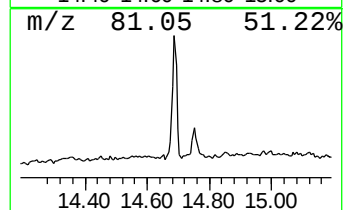
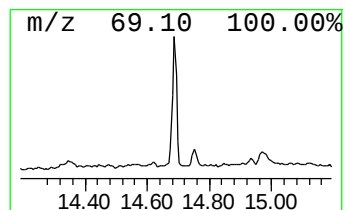
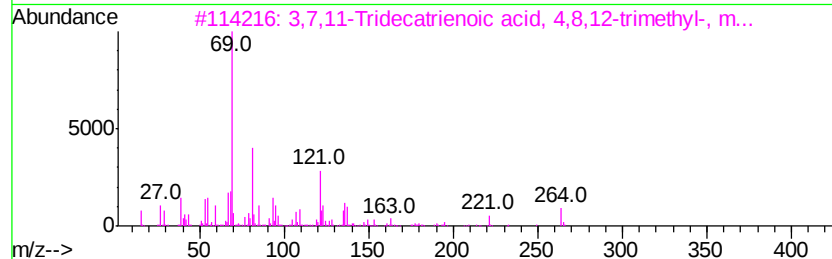
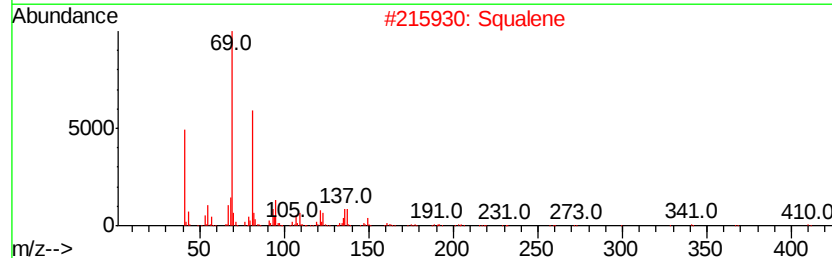
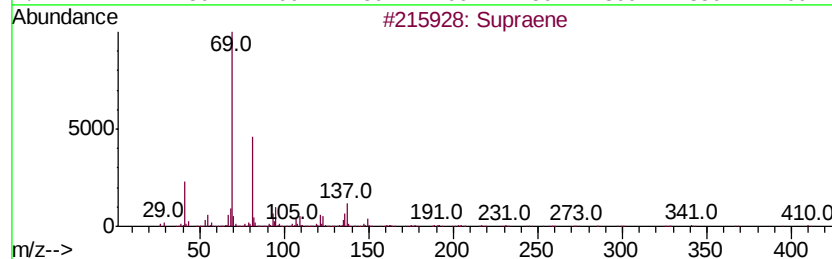
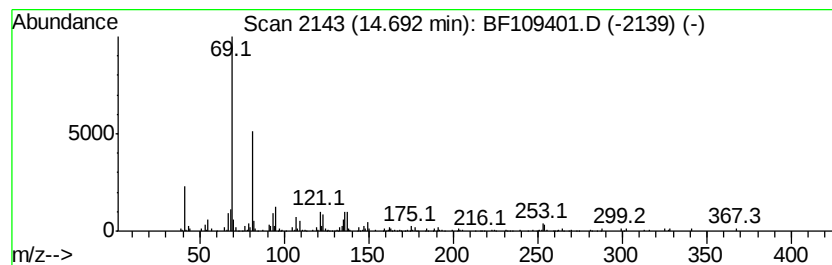
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Supraene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	8.09 ng	181320	Perylene-d12	15.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene		410	C30H50	007683-64-9	91
2	Squalene		410	C30H50	000111-02-4	91
3	3,7,11-Tridecatrienoic acid, 4,8...		264	C17H28O2	036237-70-4	64
4	Trideca-1,3,7,11-tetraene-1,1-di...		268	C18H24N2	1000159-88-9	53
5	2,3,5-Trimethyl-2,3,5-hexanetric...		203	C12H17N3	003974-79-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
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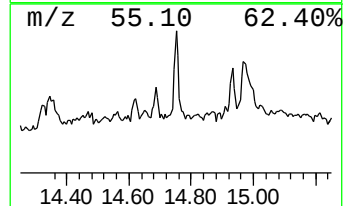
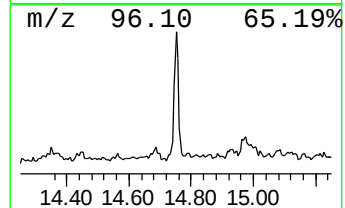
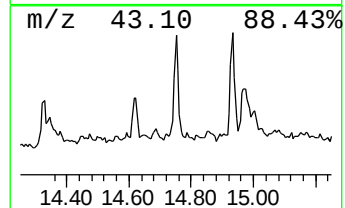
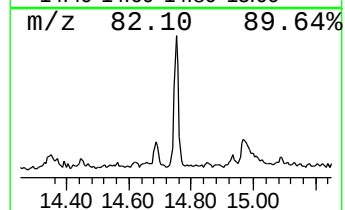
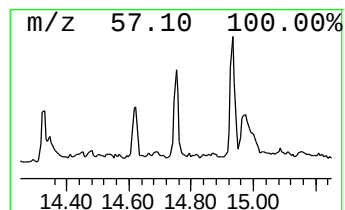
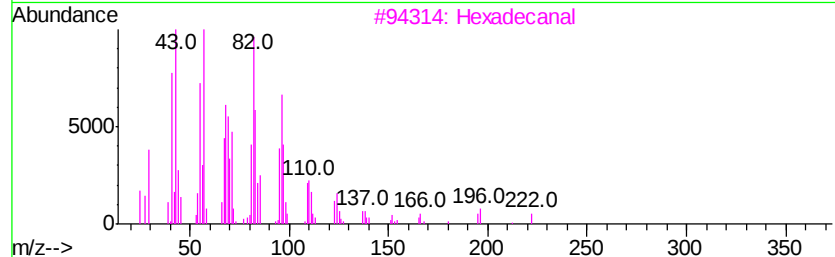
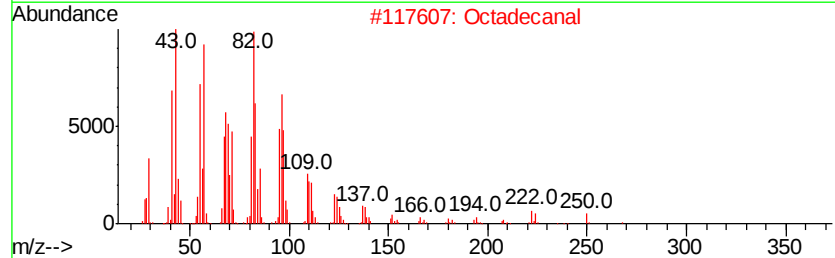
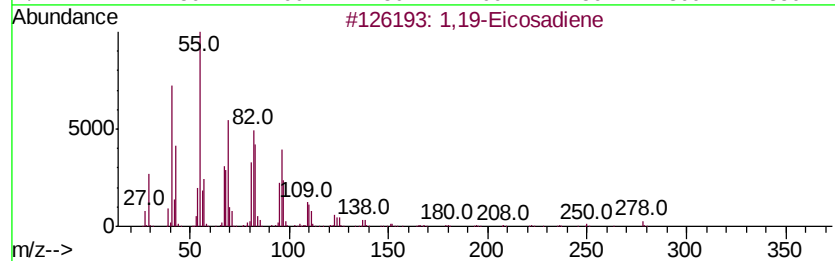
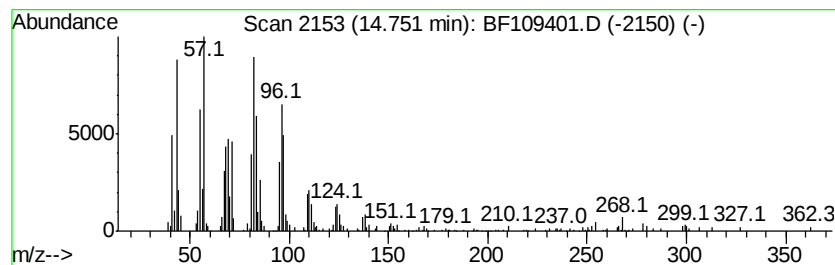
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1,19-Eicosadiene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.75	7.42 ng	166414	Perylene-d12	15.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	94
2	Octadecanal	268	C18H36O	000638-66-4	90
3	Hexadecanal	240	C16H32O	000629-80-1	90
4	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87
5	18-Nonadecen-1-ol	282	C19H38O	1000142-89-2	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
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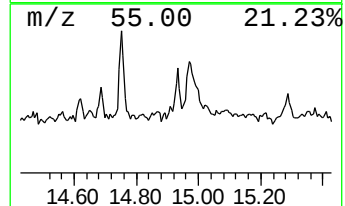
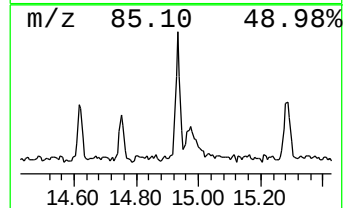
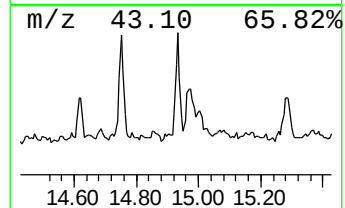
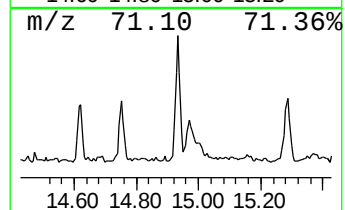
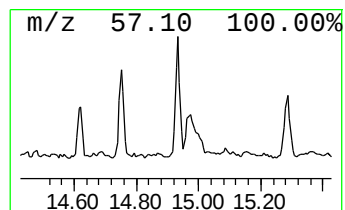
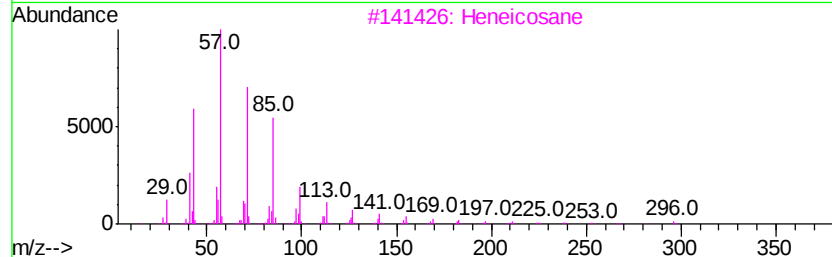
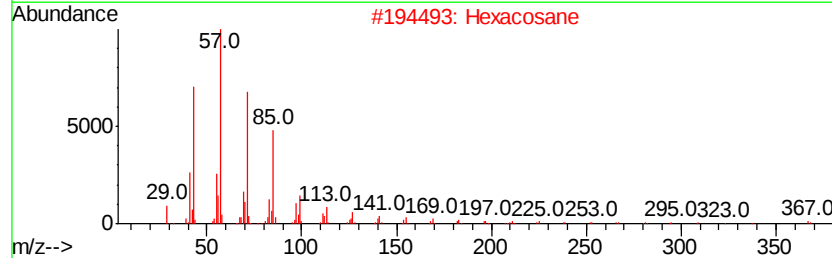
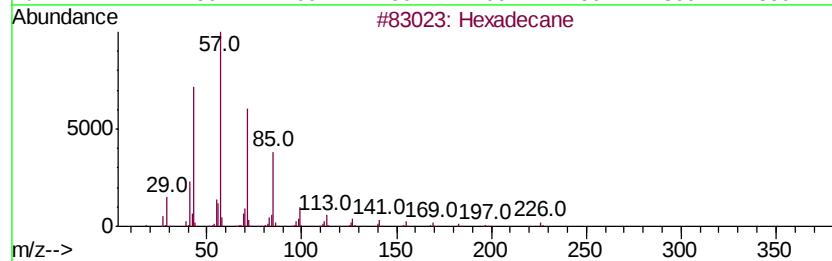
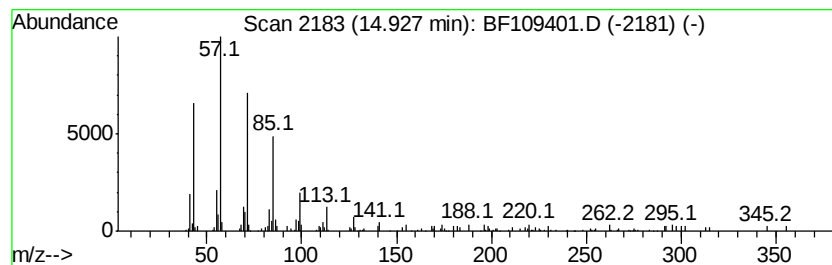
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	4.72 ng	105841	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	94
2		Hexacosane	366	C26H54	000630-01-3	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Pentacosane	352	C25H52	000629-99-2	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
Data File : BF109401.D
Acq On : 27 Sep 2018 21:28
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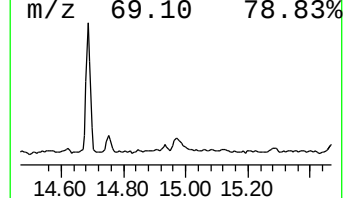
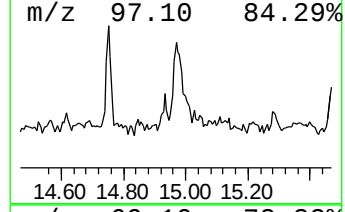
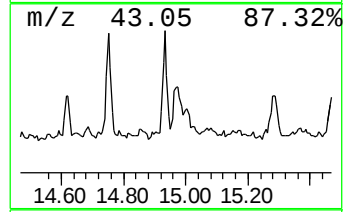
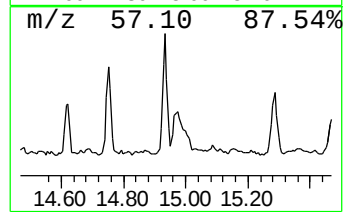
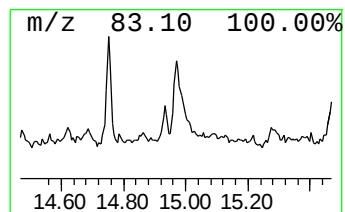
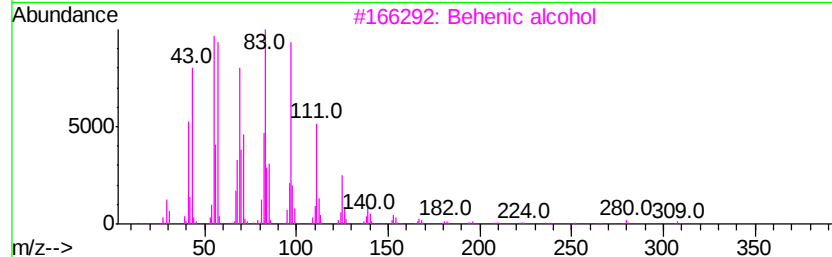
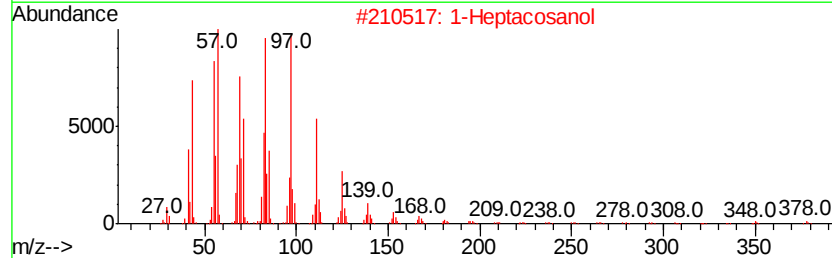
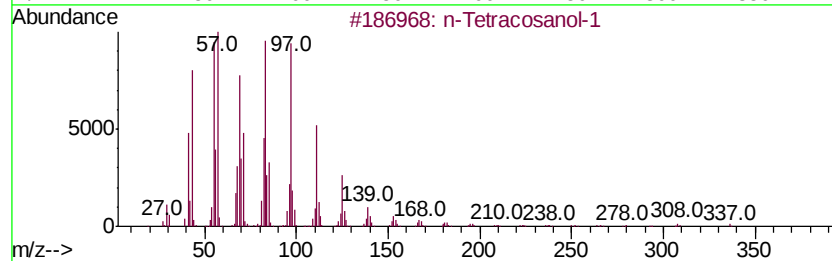
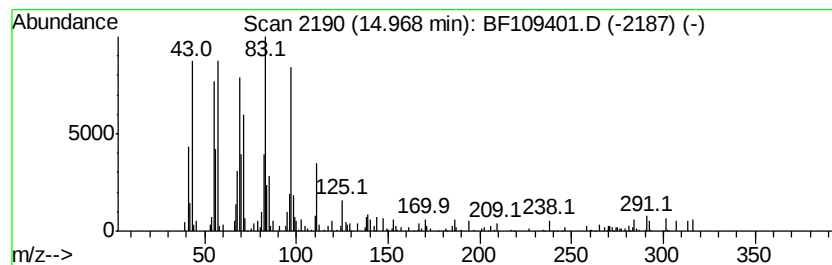
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 n-Tetracosanol-1 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	5.20 ng	116461	Perylene-d12	15.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
2			1-Heptacosanol	396	C27H56O	002004-39-9	91
3			Behenic alcohol	326	C22H46O	000661-19-8	91
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	90
5			n-Heptadecanol-1	256	C17H36O	001454-85-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
Data File : BF109401.D
Acq On : 27 Sep 2018 21:28
Operator : JU/SJ
Sample : J5077-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PARCEL-193-092418-2

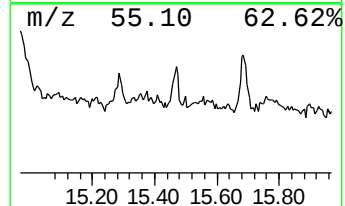
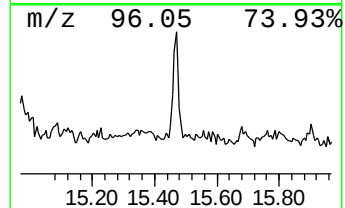
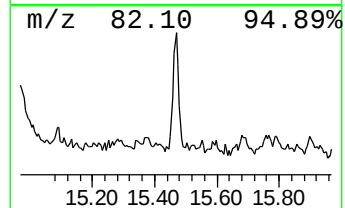
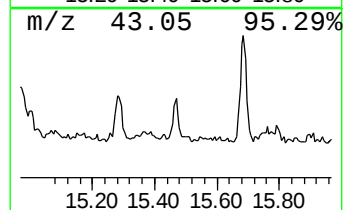
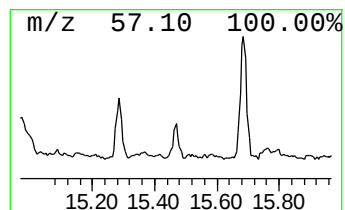
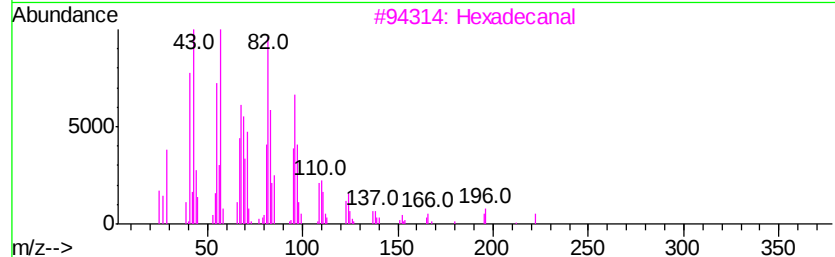
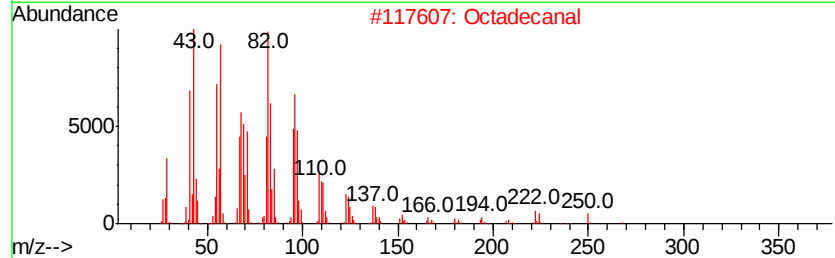
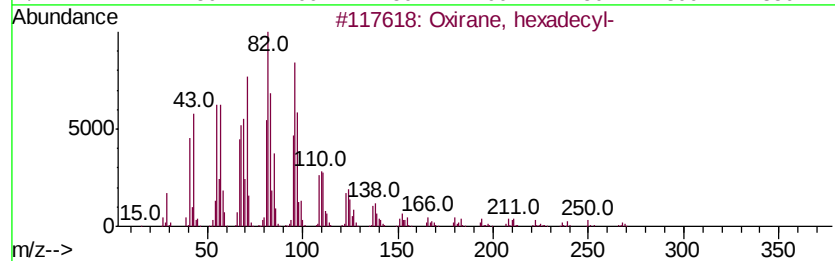
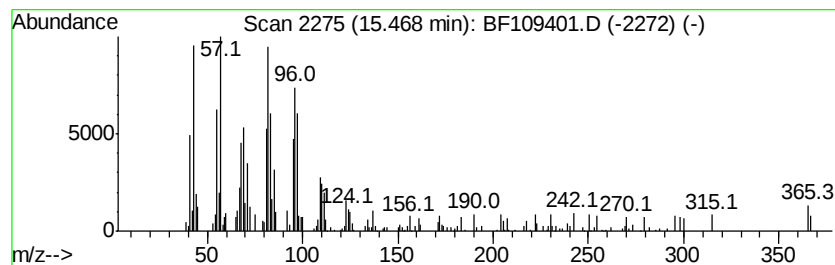
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Oxirane, hexadecyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.47	3.28 ng	73552	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
2		Octadecanal	268	C18H36O	000638-66-4	91
3		Hexadecanal	240	C16H32O	000629-80-1	91
4		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	90
5		1,15-Hexadecadiene	222	C16H30	021964-51-2	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
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Sample : J5077-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

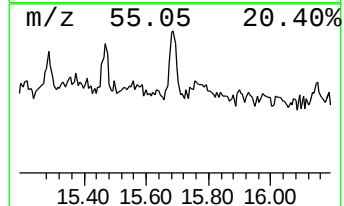
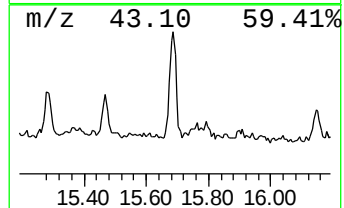
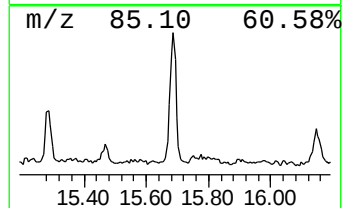
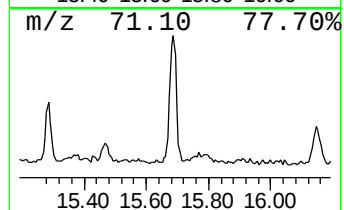
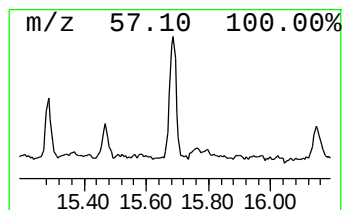
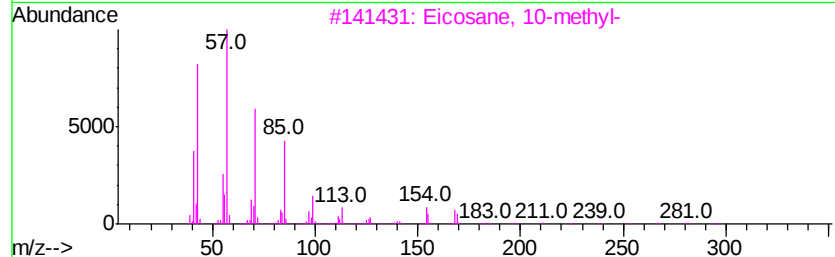
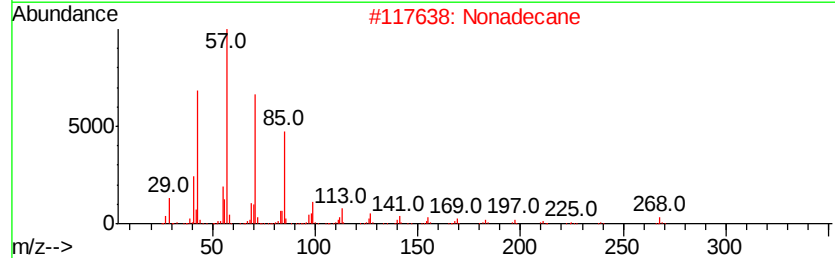
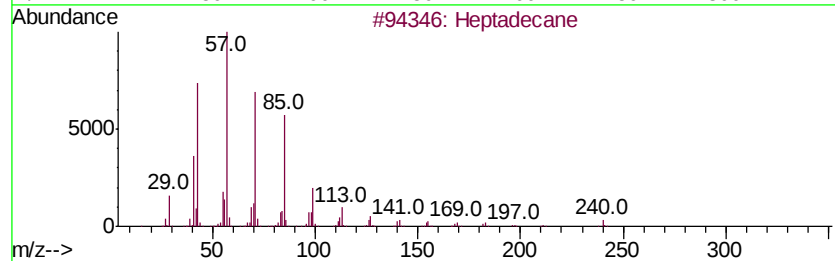
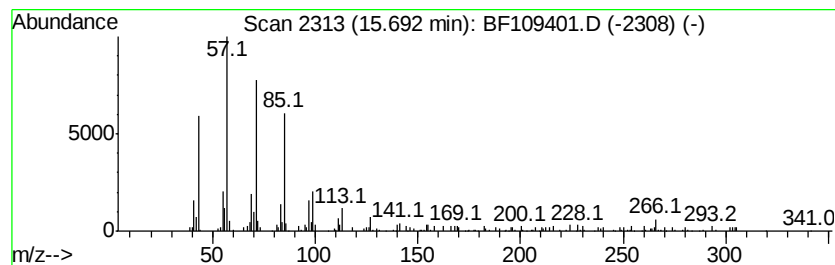
Instrument :
BNA_F
ClientSampleId :
PARCEL-193-092418-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.69	6.99 ng	156777	Perylene-d12	15.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	96
2	Nonadecane	268	C19H40	000629-92-5	93
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
4	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
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Sample : J5077-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PARCEL-193-092418-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.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
Methyl methacrylate	2.44	3.3	ng	104258	1	6.70	641469	20.0
2-Pentanone, 4-hy...	4.90	2.4	ng	75256	1	6.70	641469	20.0
unknown6.47	6.47	60.7	ng	1946200	1	6.70	641469	20.0
1-Hexadecanol	11.46	2.4	ng	80836	4	11.22	675827	20.0
n-Hexadecanoic acid	11.75	4.2	ng	140251	4	11.22	675827	20.0
Anthracene, 1-met...	11.82	2.3	ng	76781	4	11.22	675827	20.0
1-Octadecene	13.71	2.8	ng	126543	5	13.84	907958	20.0
Supraene	14.69	8.1	ng	181320	6	15.24	448278	20.0
1,19-Eicosadiene	14.75	7.4	ng	166414	6	15.24	448278	20.0
Hexadecane	14.93	4.7	ng	105841	6	15.24	448278	20.0
n-Tetracosanol-1	14.97	5.2	ng	116461	6	15.24	448278	20.0
Oxirane, hexadecyl-	15.47	3.3	ng	73552	6	15.24	448278	20.0
Heptadecane	15.69	7.0	ng	156777	6	15.24	448278	20.0