

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
 Data File : BF109400.D
 Acq On : 27 Sep 2018 21:02
 Operator : JU/SJ
 Sample : J5077-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PARCEL-193-092418-1

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	4.892	474	477	486	rBV	55021	75179	3.62%	0.335%
2	5.316	545	549	552	rBV	1829601	1651764	79.53%	7.362%
3	6.345	719	724	727	rBV	1911027	1750812	84.30%	7.803%
4	6.475	742	746	749	rBV	2086134	1772409	85.34%	7.900%
5	6.704	782	785	789	rBV	807340	658301	31.70%	2.934%
6	6.863	808	812	815	rBV	1792560	1414336	68.10%	6.304%
7	7.263	876	880	884	rBV	1163037	1123011	54.07%	5.005%
8	7.986	999	1003	1007	rBV	1028524	834284	40.17%	3.718%
9	8.686	1119	1122	1129	rBV	39396	50482	2.43%	0.225%
10	9.063	1182	1186	1189	rBV	2589890	2076877	100.00%	9.257%
11	9.457	1249	1253	1256	rBV	818330	610997	29.42%	2.723%
12	9.733	1297	1300	1305	rBV	959124	884964	42.61%	3.944%
13	10.521	1430	1434	1440	rBV	1101107	1017667	49.00%	4.536%
14	11.216	1548	1552	1554	rBV	821246	755293	36.37%	3.366%
15	11.233	1554	1555	1558	rVB	97922	64400	3.10%	0.287%
16	11.716	1634	1637	1639	rBV	32686	26382	1.27%	0.118%
17	11.751	1639	1643	1648	rVV2	137055	148955	7.17%	0.664%
18	11.821	1652	1655	1658	rBV	26653	27197	1.31%	0.121%
19	12.263	1726	1730	1732	rBV	24567	25138	1.21%	0.112%
20	12.415	1753	1756	1760	rBV	161692	151030	7.27%	0.673%
21	12.533	1774	1776	1781	rBV3	23510	28399	1.37%	0.127%
22	12.645	1790	1795	1798	rBV	167103	160084	7.71%	0.713%
23	12.792	1816	1820	1824	rVB	1673052	1492686	71.87%	6.653%
24	12.974	1849	1851	1855	rVB	28151	25068	1.21%	0.112%
25	13.039	1856	1862	1865	rBV3	50657	68523	3.30%	0.305%
26	13.380	1917	1920	1926	rVB6	22612	27409	1.32%	0.122%
27	13.492	1938	1939	1945	rVB3	30743	27121	1.31%	0.121%
28	13.586	1952	1955	1957	rBV2	19904	25978	1.25%	0.116%
29	13.709	1973	1976	1987	rVB2	160209	257068	12.38%	1.146%
30	13.839	1993	1998	2001	rBV	869173	874155	42.09%	3.896%
31	13.862	2001	2002	2005	rVB	92312	51347	2.47%	0.229%
32	14.021	2027	2029	2032	rBV3	29416	27305	1.31%	0.122%
33	14.139	2047	2049	2054	rVB5	34811	37822	1.82%	0.169%
34	14.345	2077	2084	2093	rBV4	170554	403734	19.44%	1.799%

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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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.451	2099	2102	2104	rBV4	25362	28765	1.39%	0.128%
36	14.474	2104	2106	2111	rVB	82091	79590	3.83%	0.355%
37	14.621	2128	2131	2133	rBV	51962	50523	2.43%	0.225%
38	14.686	2139	2142	2146	rVB	152814	147908	7.12%	0.659%
39	14.751	2150	2153	2158	rVB	467852	431970	20.80%	1.925%
40	14.845	2166	2169	2178	rVB	80522	137386	6.62%	0.612%
41	14.933	2180	2184	2187	rVV	215848	237952	11.46%	1.061%
42	14.968	2187	2190	2204	rVB2	215581	460576	22.18%	2.053%
43	15.121	2214	2216	2220	rVB	56383	56490	2.72%	0.252%
44	15.180	2223	2226	2230	rVV	54808	57088	2.75%	0.254%
45	15.239	2232	2236	2240	rVV	447045	515528	24.82%	2.298%
46	15.286	2241	2244	2247	rVB	68194	84804	4.08%	0.378%
47	15.468	2271	2275	2282	rVB2	149954	201990	9.73%	0.900%
48	15.686	2307	2312	2317	rBV	269280	360030	17.34%	1.605%
49	16.415	2431	2436	2441	rVB2	79221	128957	6.21%	0.575%
50	16.686	2478	2482	2490	rVB2	90251	183325	8.83%	0.817%
51	17.645	2641	2645	2651	rVB5	49505	97476	4.69%	0.434%
52	17.886	2678	2686	2708	rVB8	127821	550018	26.48%	2.451%

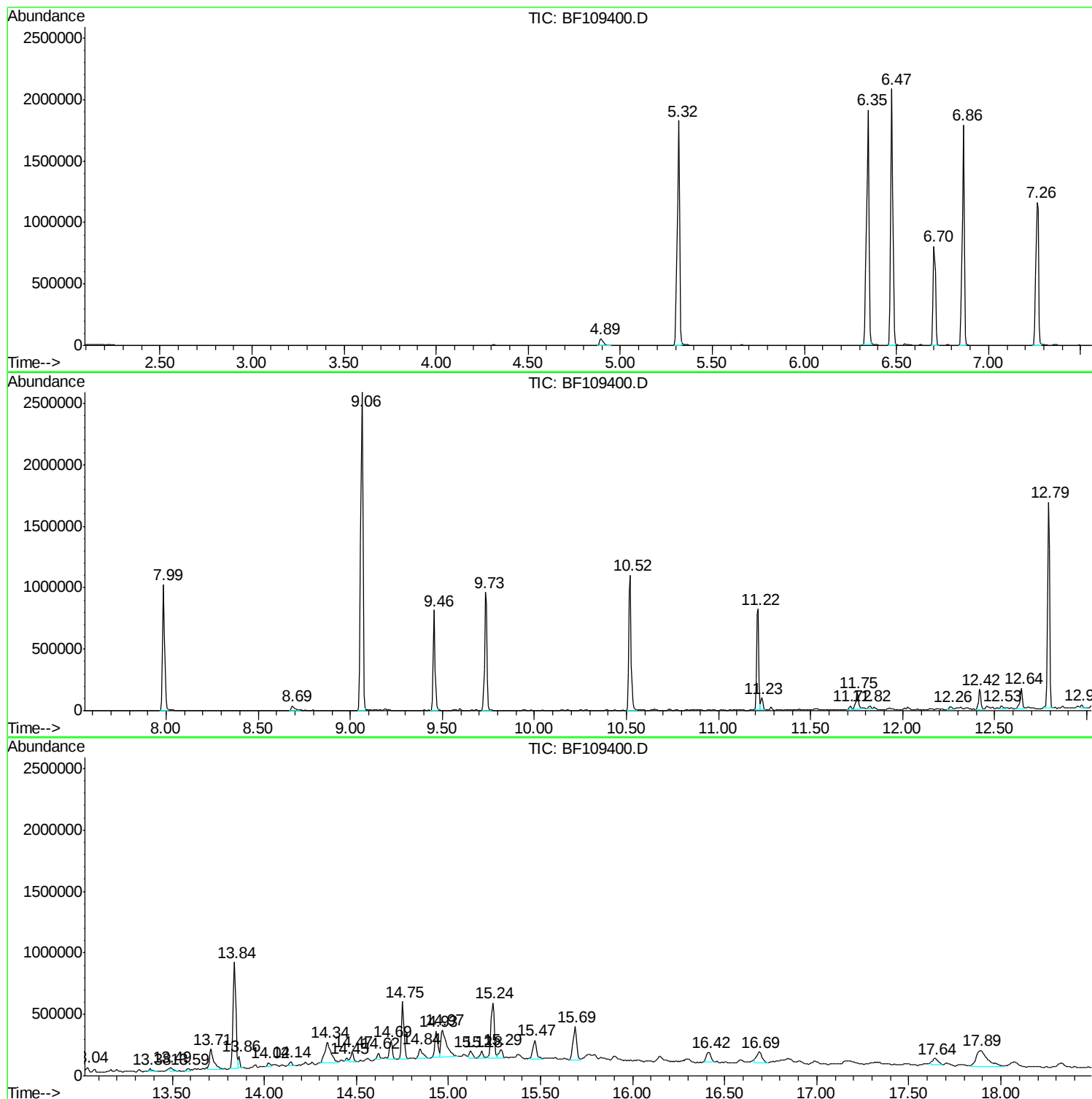
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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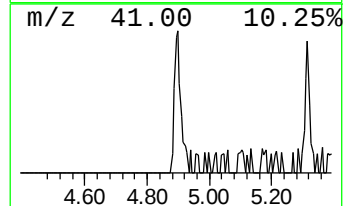
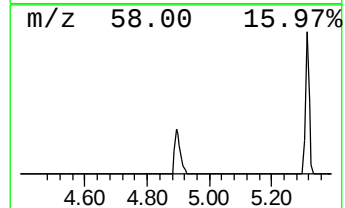
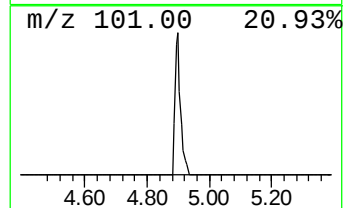
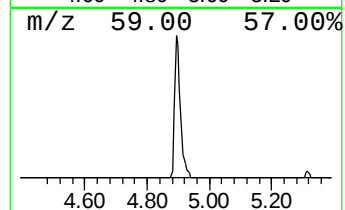
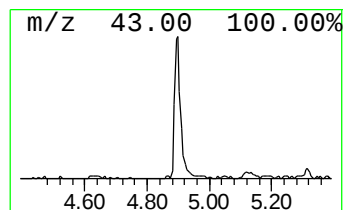
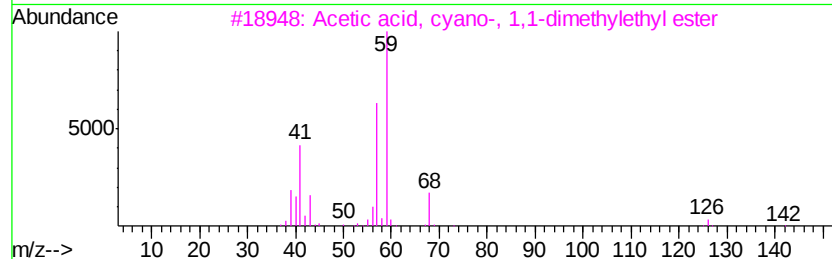
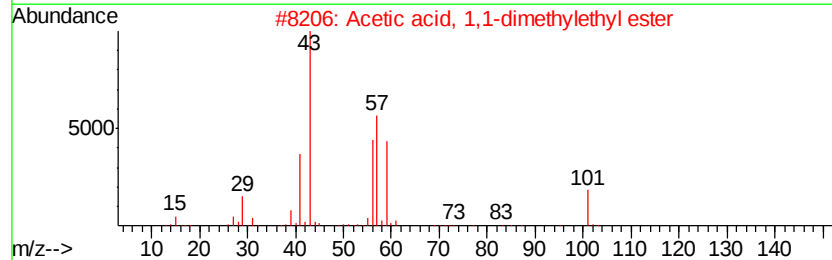
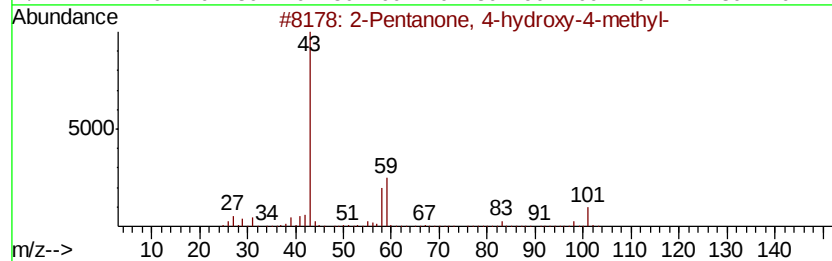
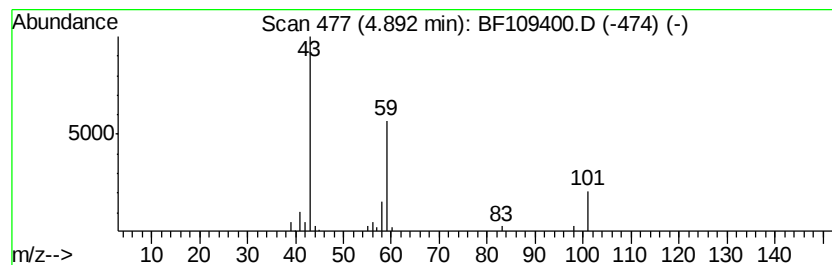
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	2.28 ng	75179	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			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	9



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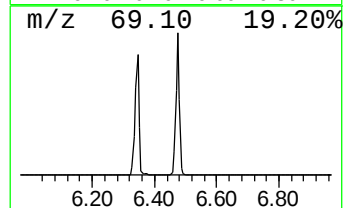
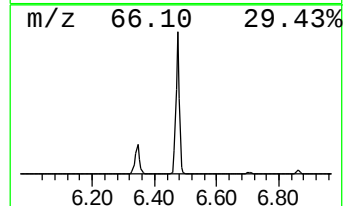
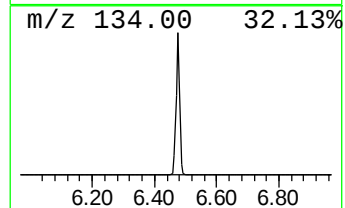
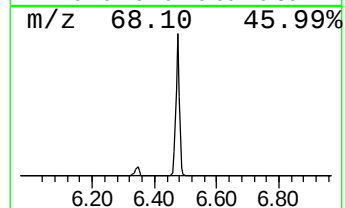
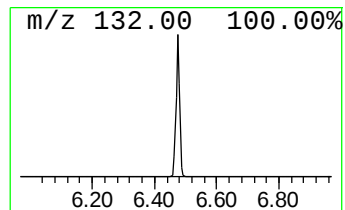
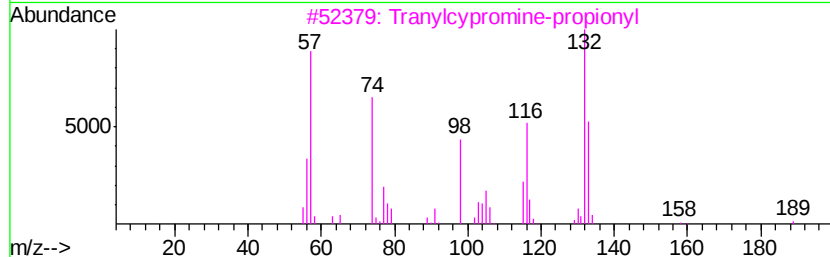
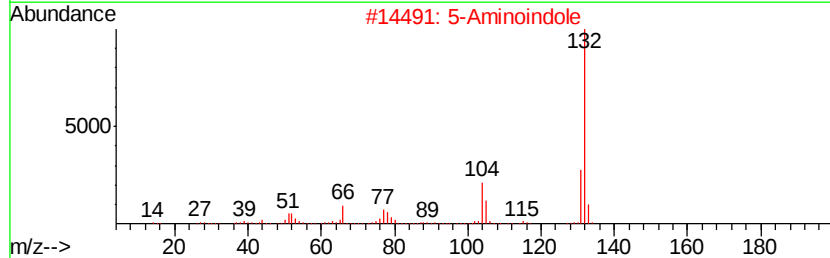
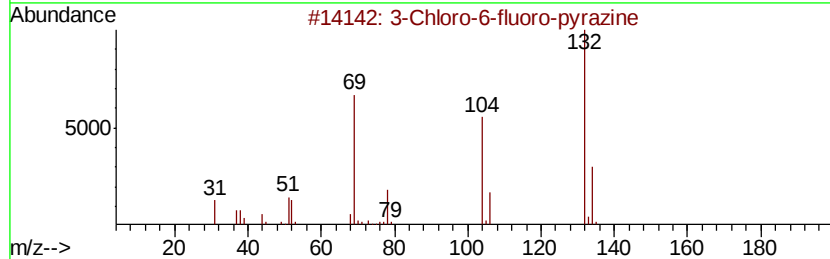
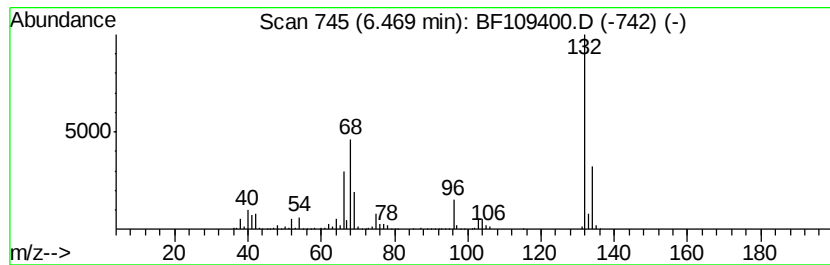
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Peak Number 2 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	53.85 ng	1772410	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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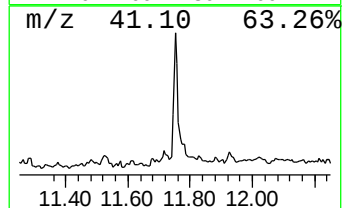
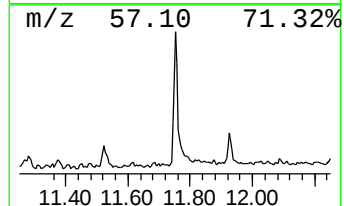
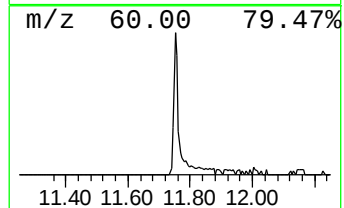
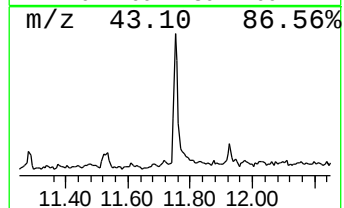
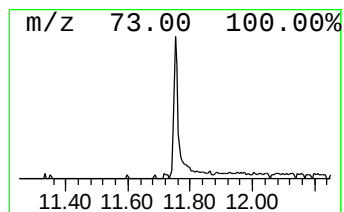
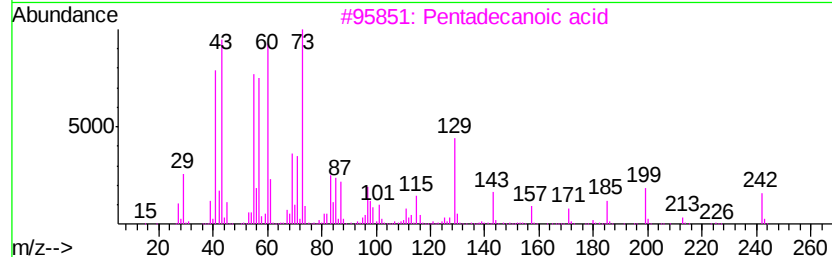
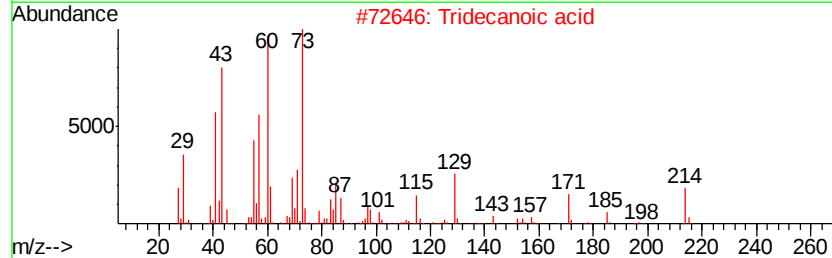
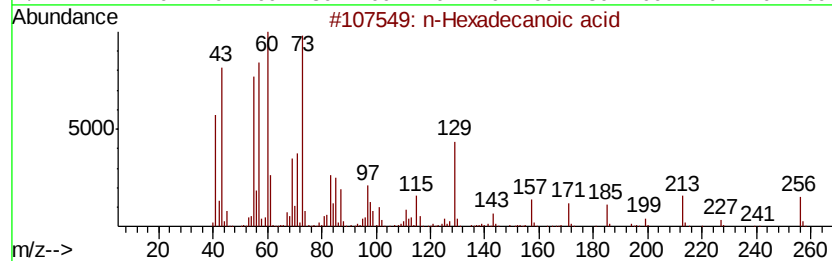
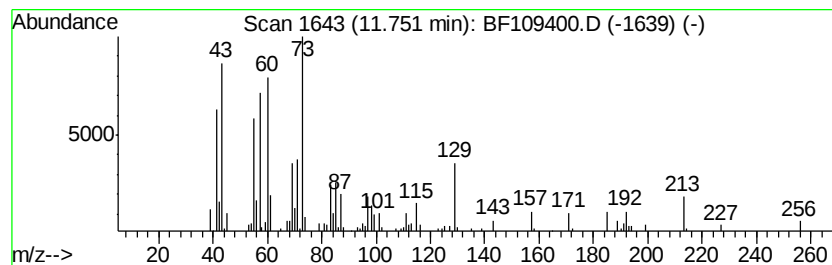
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	3.94 ng	148955	Phenanthrene-d10	11.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5		Dodecanoic acid	200	C12H24O2	000143-07-7	72



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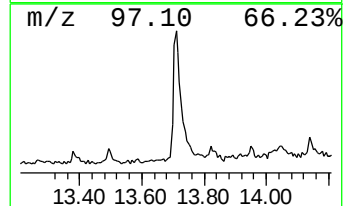
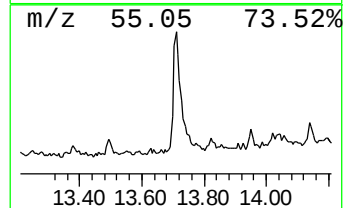
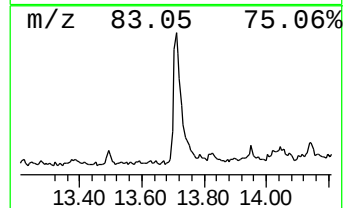
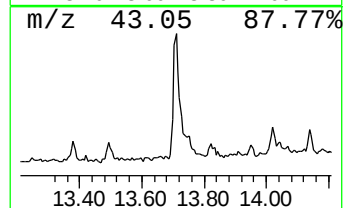
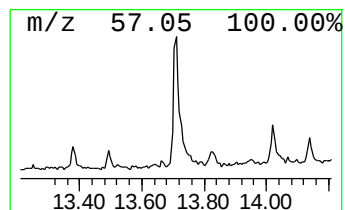
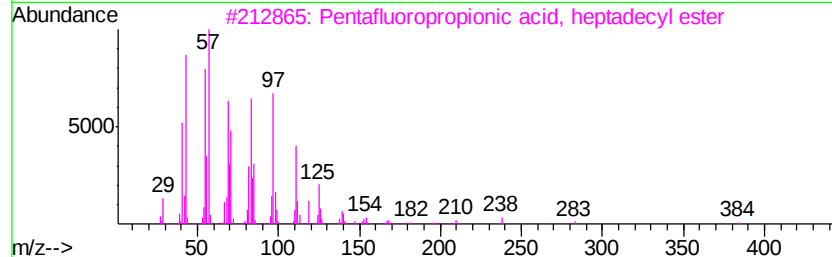
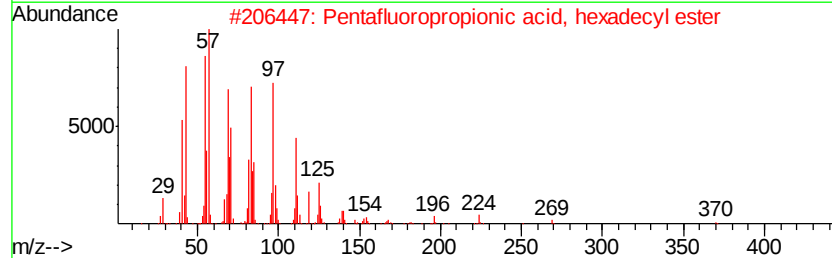
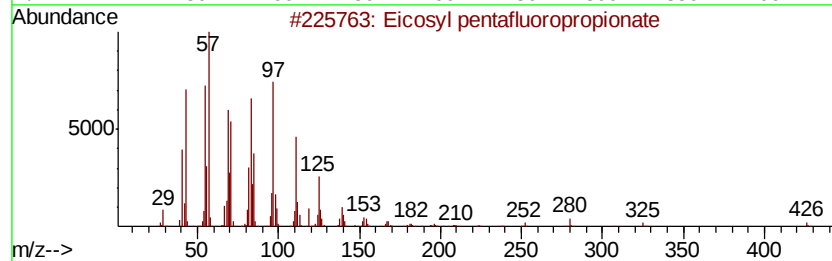
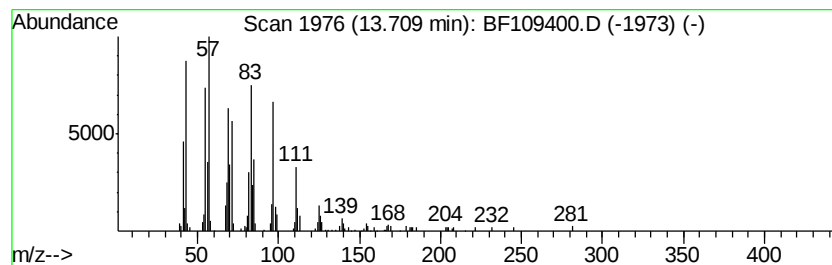
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Eicosyl pentafluoropropionate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	5.88 ng	257068	Chrysene-d12	13.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	91	
2	Pentafluoropropionic acid, hexadecyl ester	388	C19H33F5O2	006222-07-7	91	
3	Pentafluoropropionic acid, heptadecyl ester	402	C20H35F5O2	959218-78-1	91	
4	Trichloroacetic acid, hexadecyl ester	386	C18H33Cl3O2	074339-54-1	91	
5	Heptafluorobutyric acid, pentadecyl ester	424	C19H31F7O2	959261-23-5	91	



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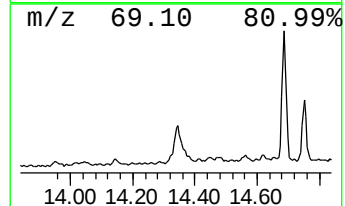
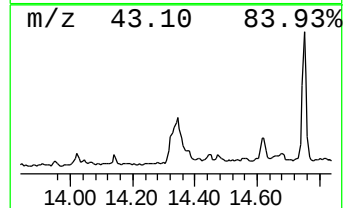
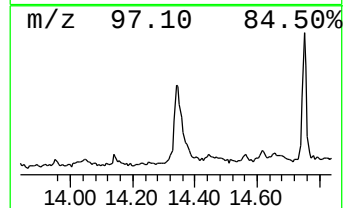
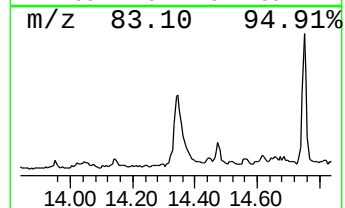
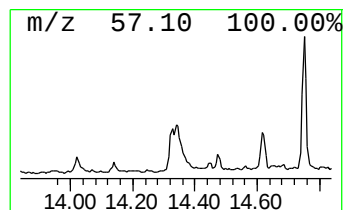
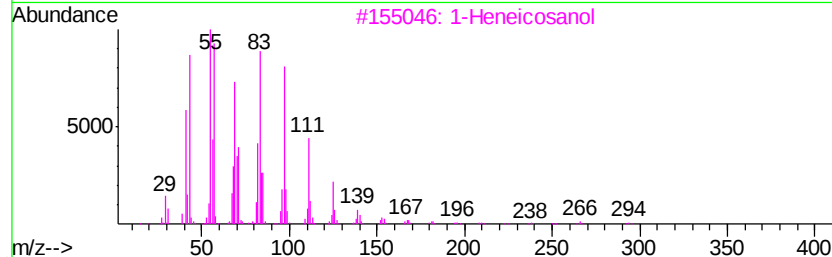
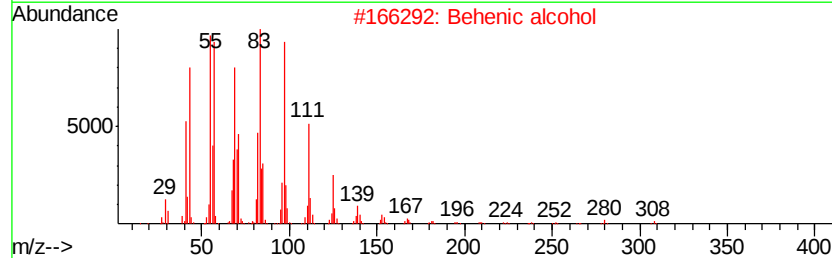
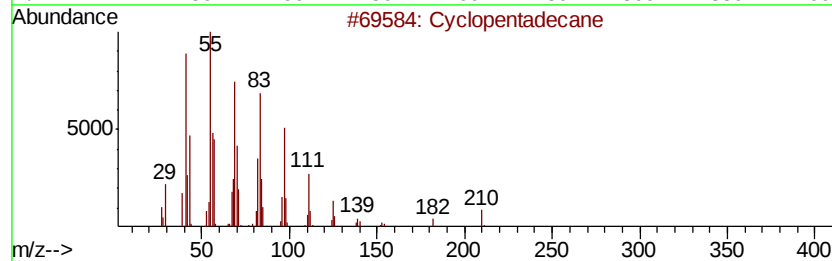
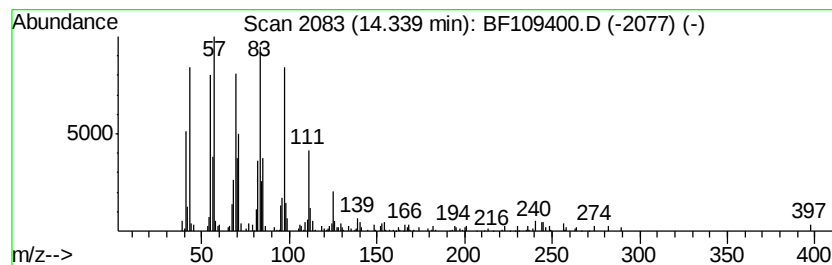
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Cyclopentadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.34	9.24 ng	403734	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	98
2		Behenic alcohol	326	C22H46O	000661-19-8	95
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	93
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
Operator : JU/SJ
Sample : J5077-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

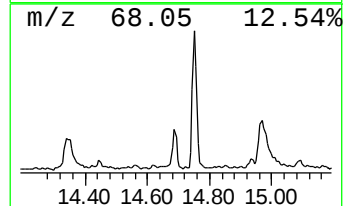
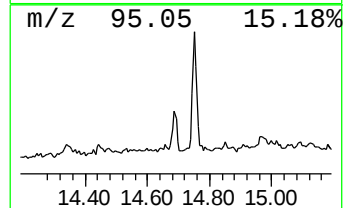
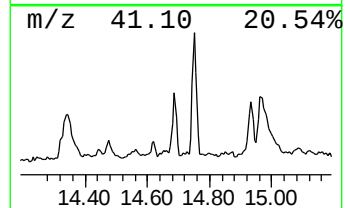
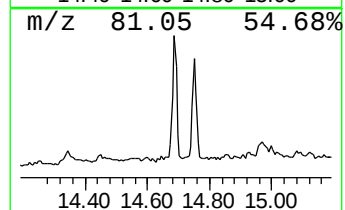
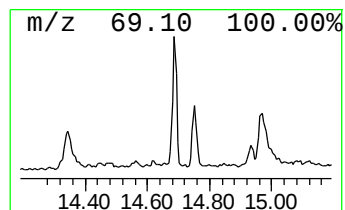
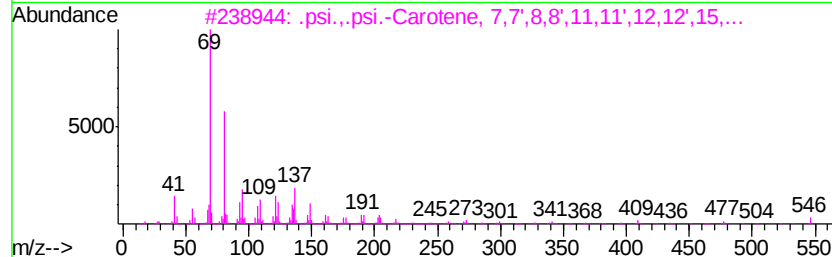
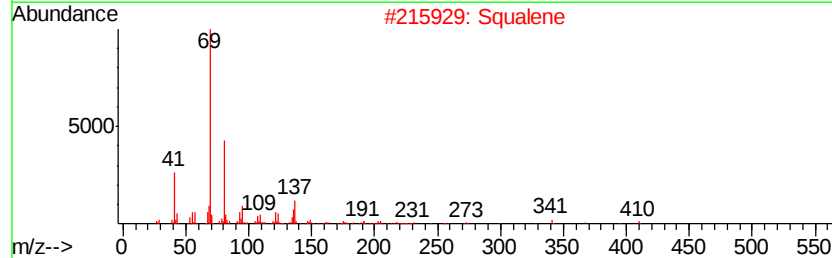
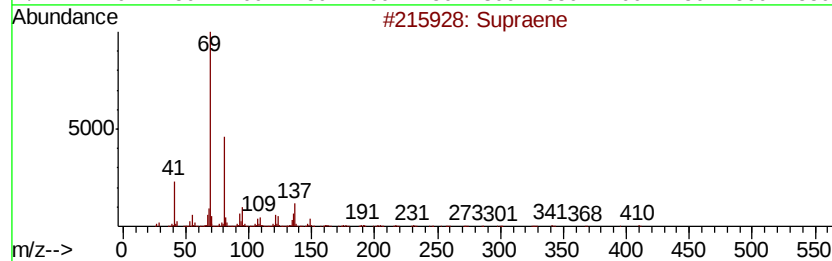
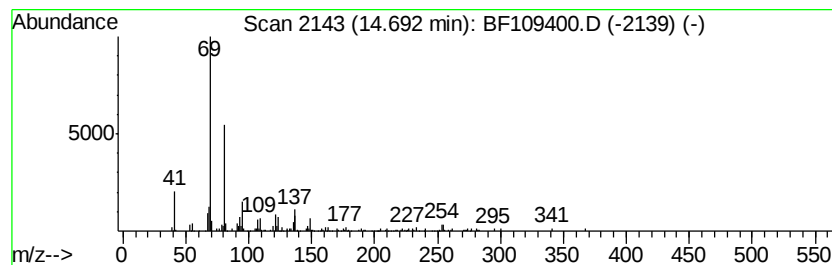
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Supraene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	5.74 ng	147908	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 80
3	.psi.,.psi.-Carotene, 7,7',8,8',...		547 C40H66	000502-62-5 78
4	3,7,11-Tridecatrienitrile, 4,8...		231 C16H25N	006006-01-5 72
5	2,6,10,14-Hexadecatetraenoic aci...		332 C22H36O2	024035-35-6 64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
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Operator : JU/SJ
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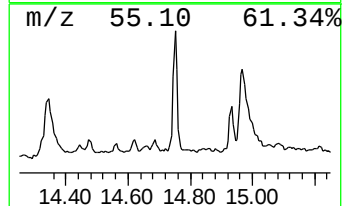
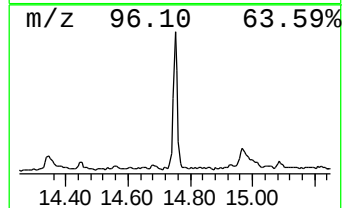
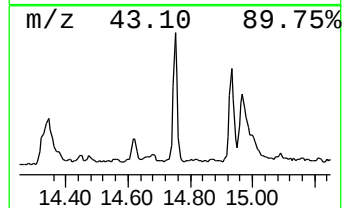
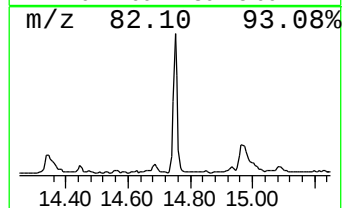
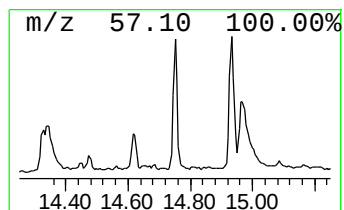
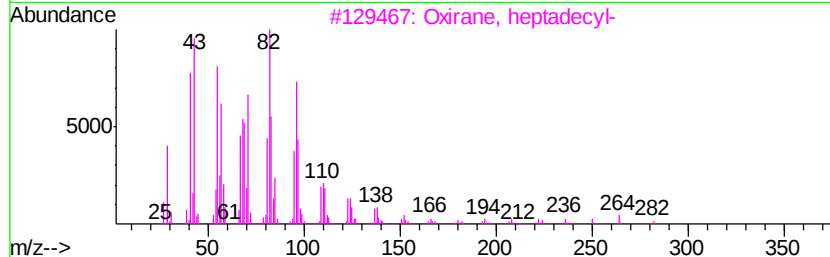
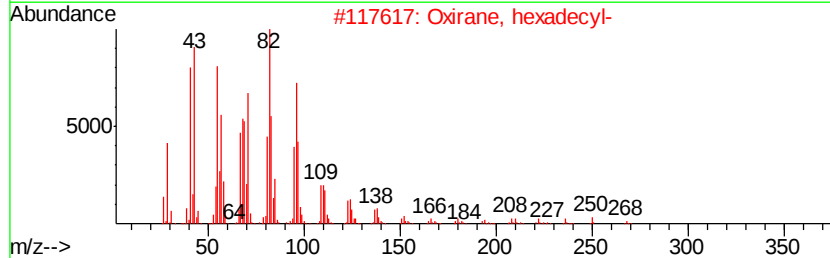
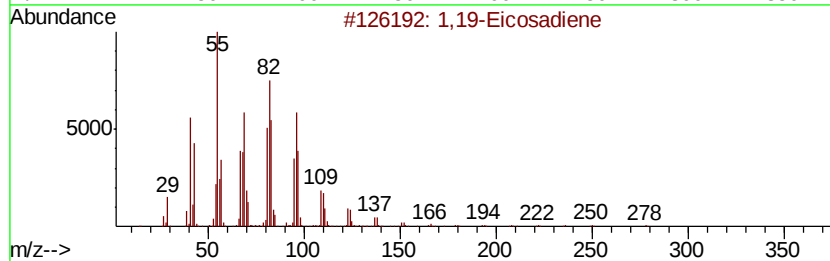
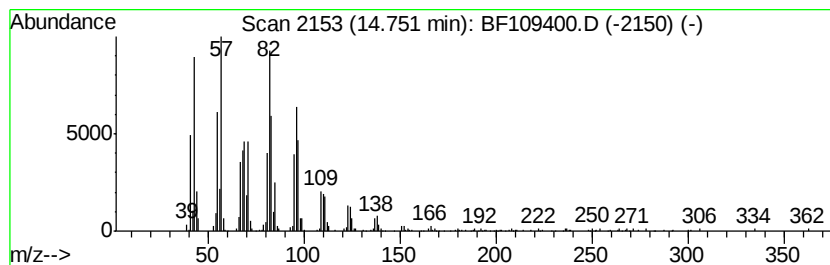
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1,19-Eicosadiene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	16.76 ng	431970	Perylene-d12	15.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,19-Eicosadiene	278	C20H38	014811-95-1	95
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	93
3		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4		Octadecanal	268	C18H36O	000638-66-4	91
5		Z-2-Octadecen-1-ol	268	C18H36O	1000131-11-0	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
Operator : JU/SJ
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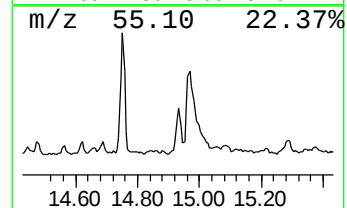
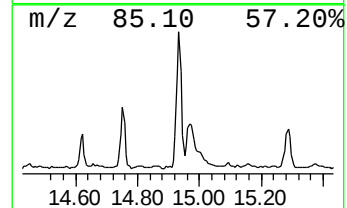
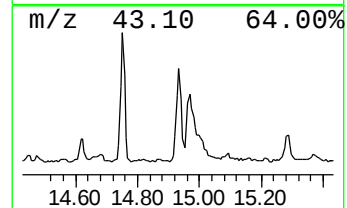
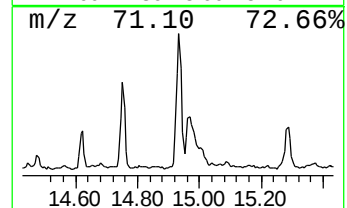
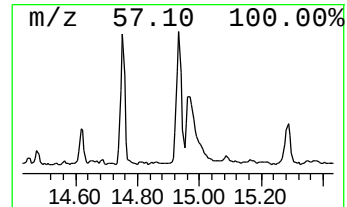
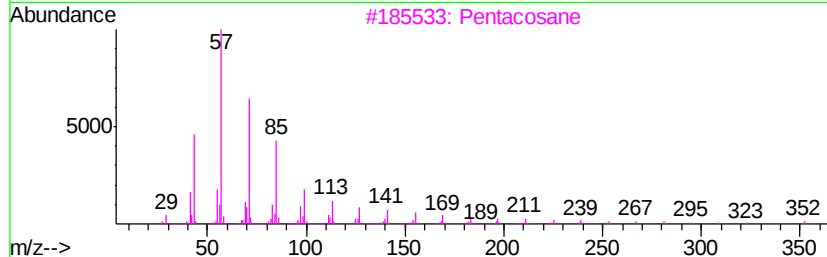
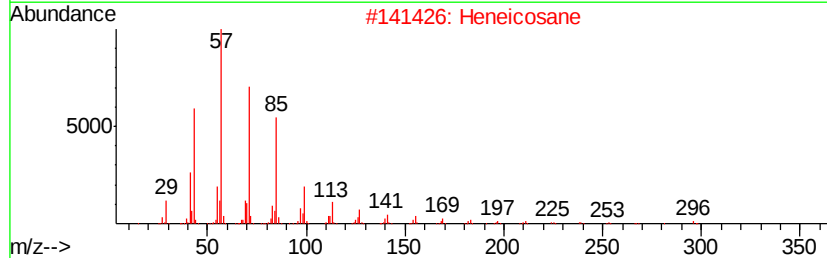
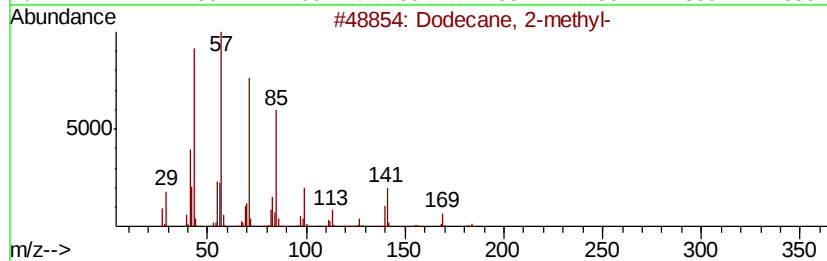
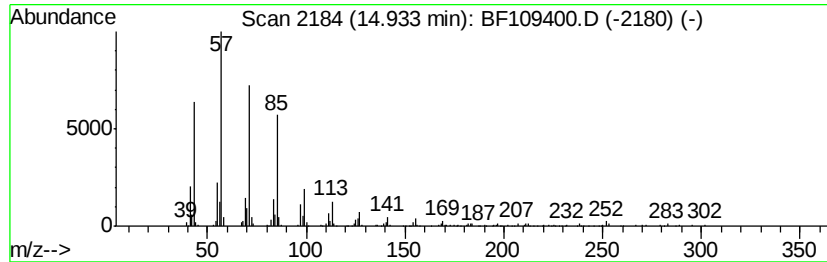
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Dodecane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	9.23 ng	237952	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
2		Heneicosane	296	C21H44	000629-94-7	90
3		Pentacosane	352	C25H52	000629-99-2	87
4		Hexadecane, 2-methyl-	240	C17H36	001560-92-5	87
5		Eicosane	282	C20H42	000112-95-8	86



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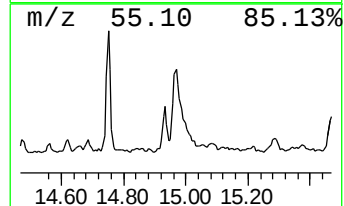
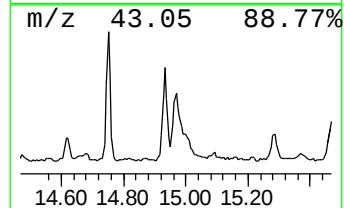
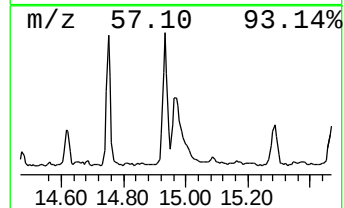
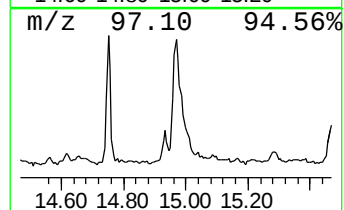
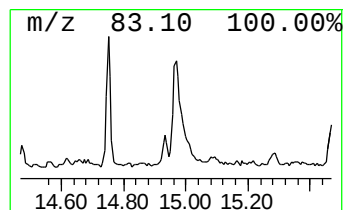
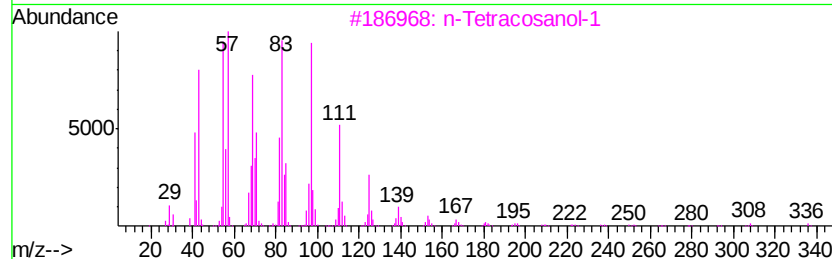
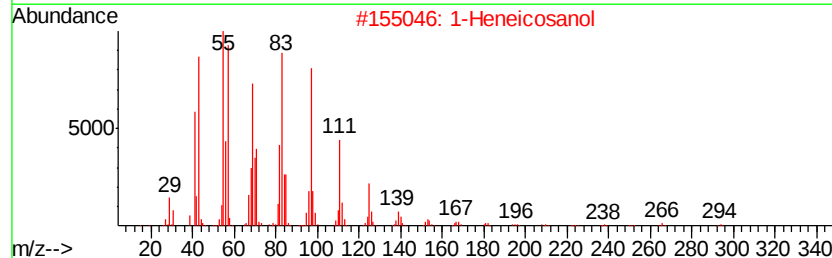
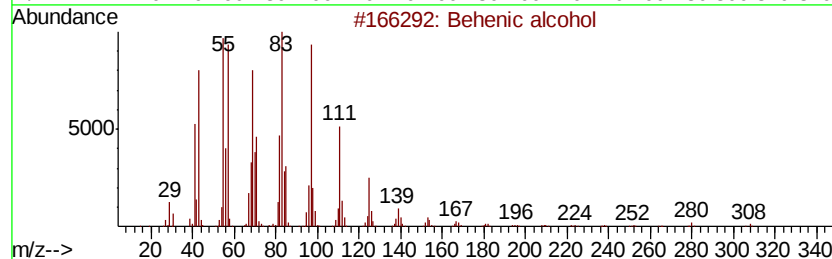
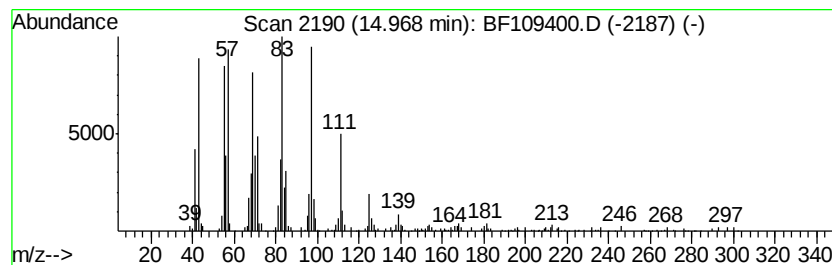
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Behenic alcohol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	17.87 ng	460576	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Behenic alcohol	326	C22H46O	000661-19-8	95
2		1-Heneicosanol	312	C21H44O	015594-90-8	94
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	92
4		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	76
5		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
Operator : JU/SJ
Sample : J5077-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PARCEL-193-092418-1

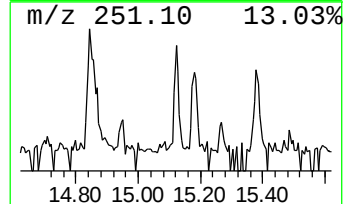
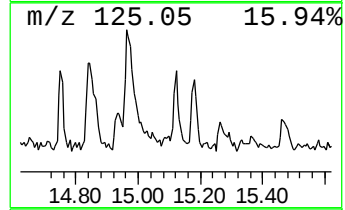
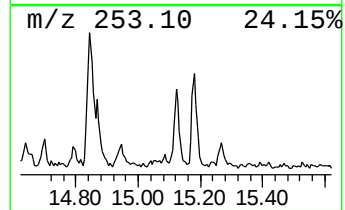
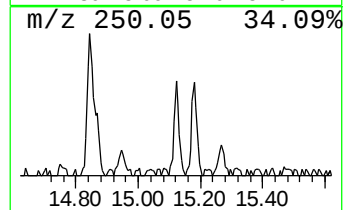
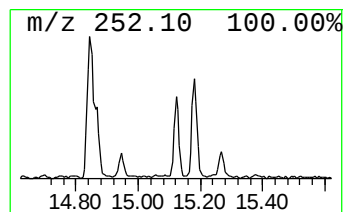
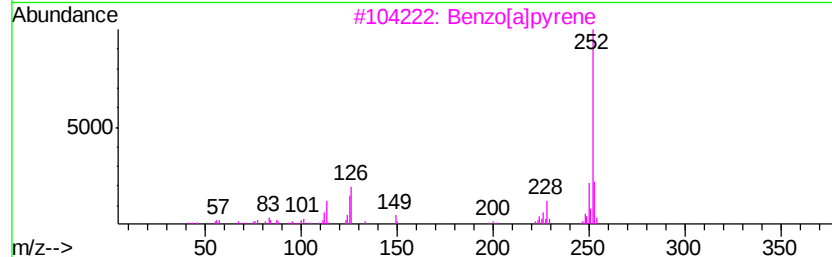
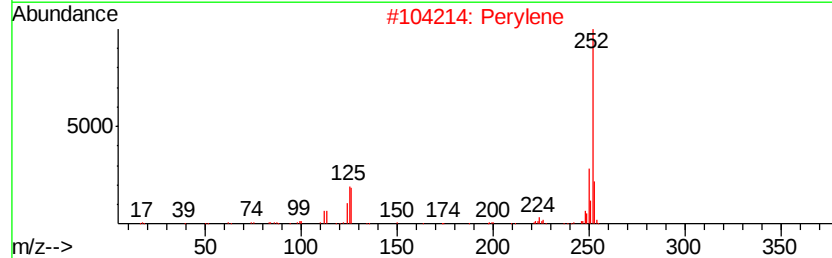
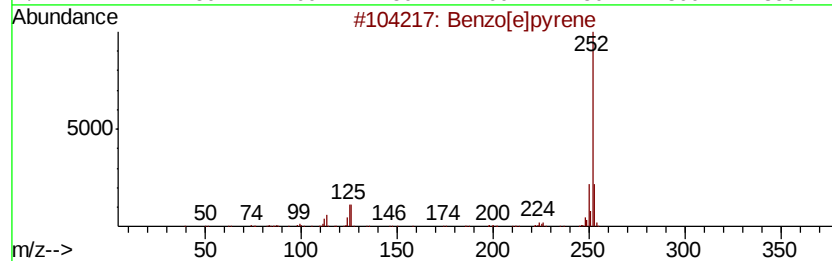
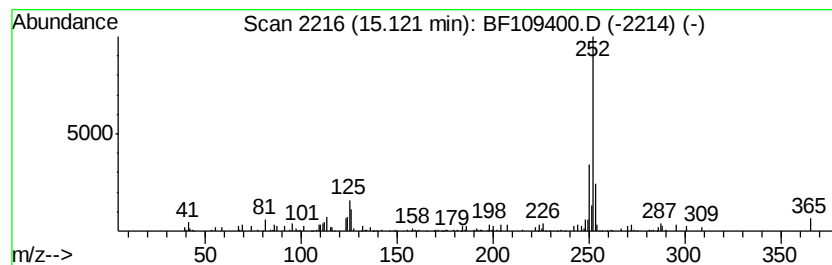
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	2.19 ng	56490	Perylene-d12	15.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	96
2			Perylene	252	C20H12	000198-55-0	90
3			Benzo[a]pyrene	252	C20H12	000050-32-8	81
4			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	76
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	70



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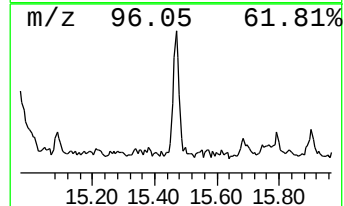
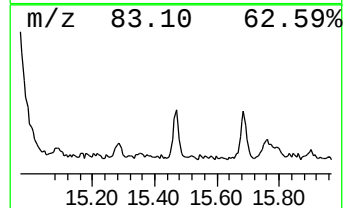
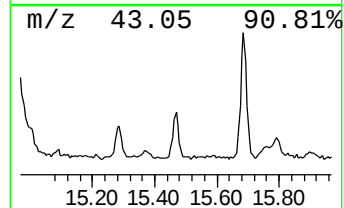
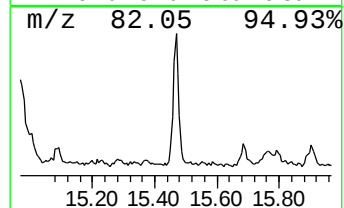
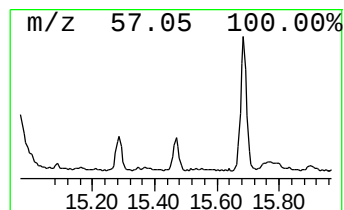
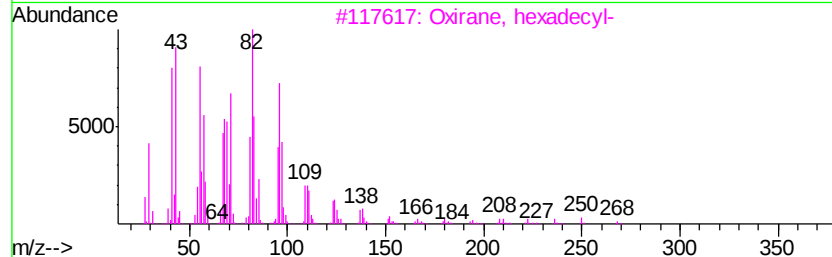
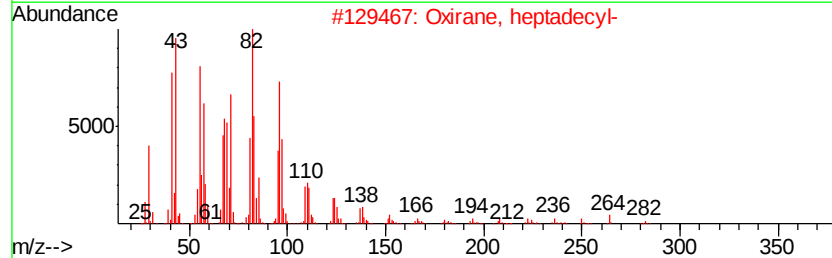
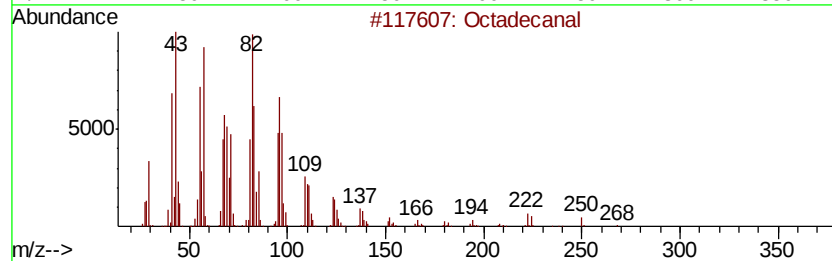
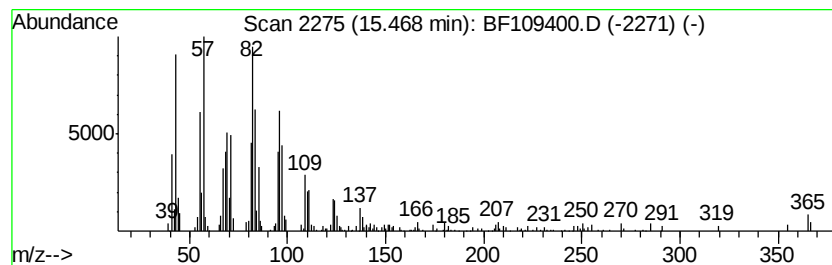
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecanal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.47	7.84 ng	201990	Perylene-d12		15.24	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanal		268	C18H36O	000638-66-4	91
2	Oxirane, heptadecyl-		282	C19H38O	067860-04-2	91
3	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	90
4	Hexadecanal		240	C16H32O	000629-80-1	90
5	13-Octadecenal		266	C18H34O	056554-90-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
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Sample : J5077-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PARCEL-193-092418-1

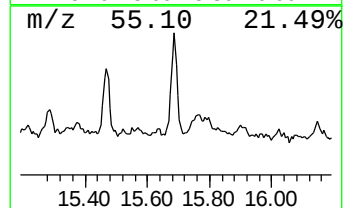
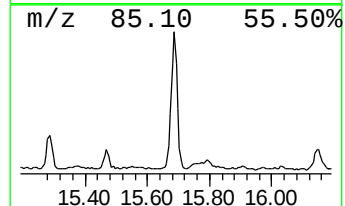
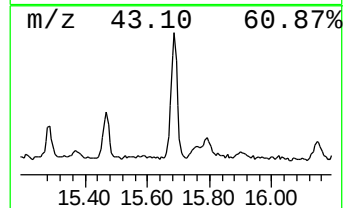
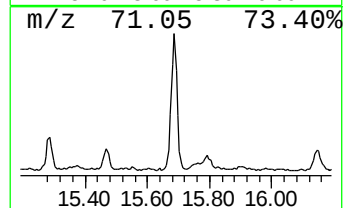
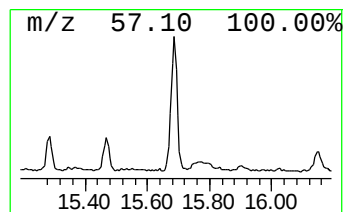
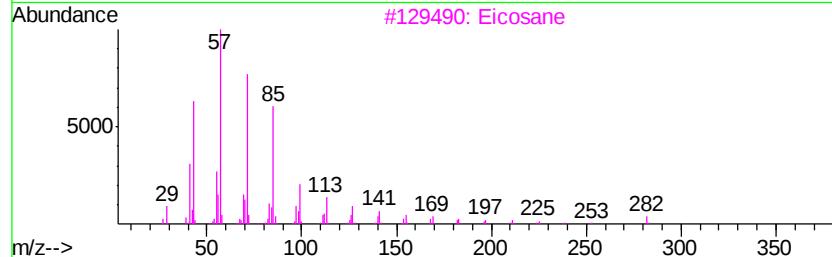
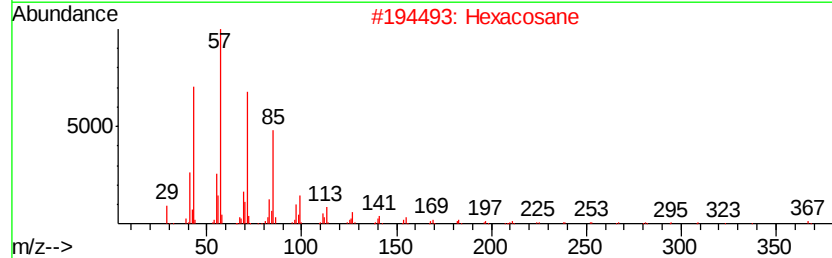
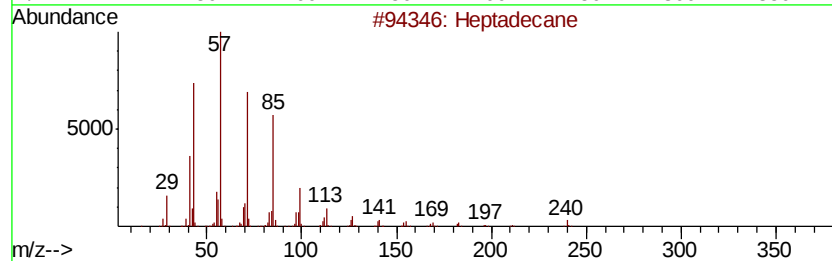
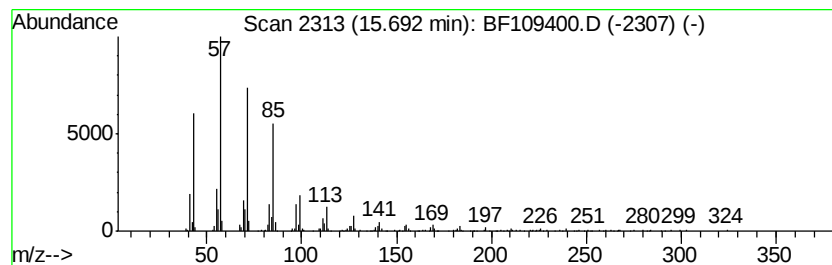
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	13.97 ng	360030	Perylene-d12	15.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	94
2	Hexacosane		366	C26H54	000630-01-3	91
3	Eicosane		282	C20H42	000112-95-8	91
4	Hexadecane		226	C16H34	000544-76-3	91
5	Heneicosane		296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
Operator : JU/SJ
Sample : J5077-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

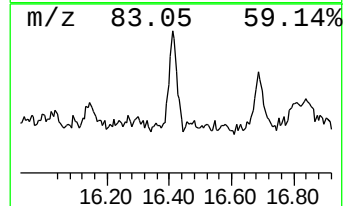
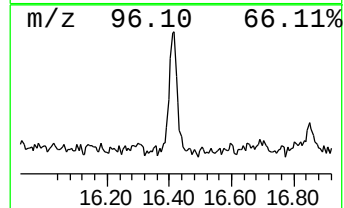
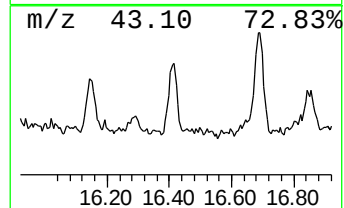
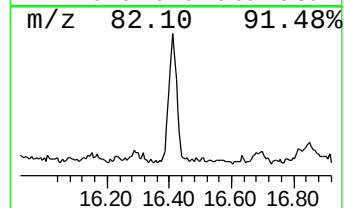
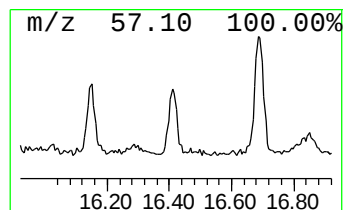
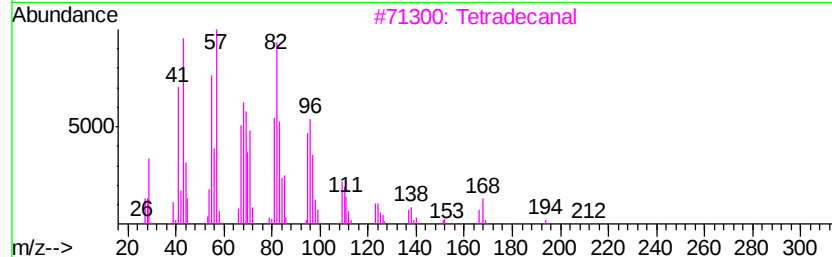
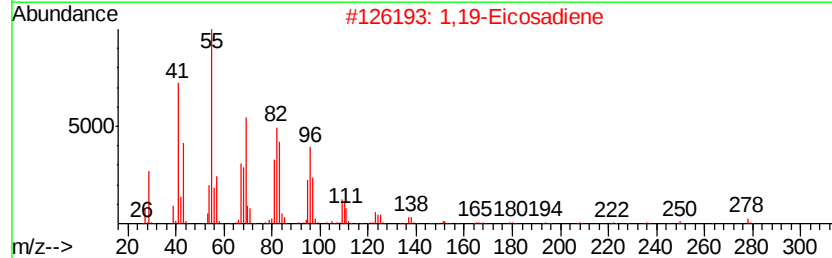
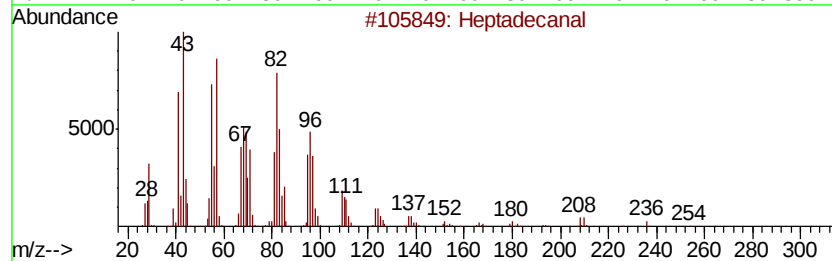
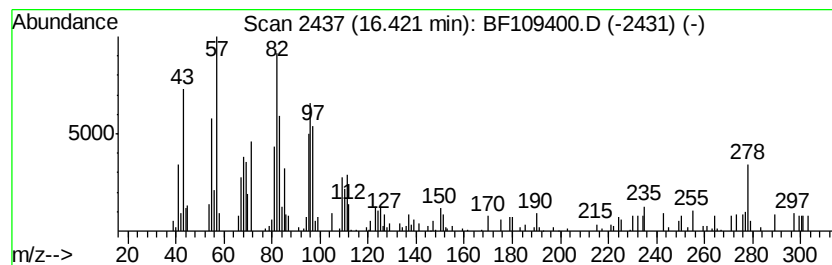
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ClientSampleId :
PARCEL-193-092418-1

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 14 Heptadecanal Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.42	5.00 ng	128957	Perylene-d12	15.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecanal	254	C17H34O	1000376-70-0	86
2	1,19-Eicosadiene	278	C20H38	014811-95-1	83
3	Tetradecanal	212	C14H28O	000124-25-4	80
4	1,15-Pentadecanediol	244	C15H32O2	014722-40-8	72
5	1,15-Hexadecadiene	222	C16H30	021964-51-2	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
Operator : JU/SJ
Sample : J5077-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PARCEL-193-092418-1

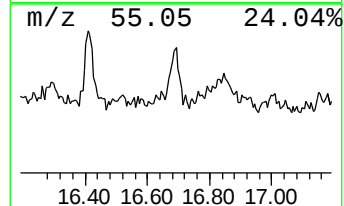
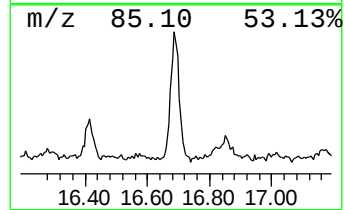
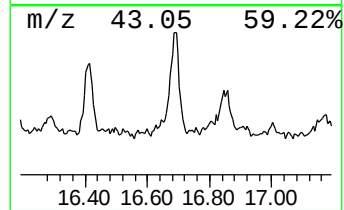
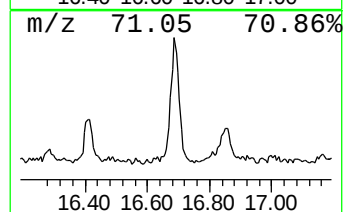
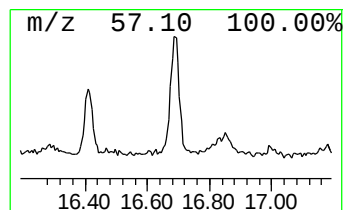
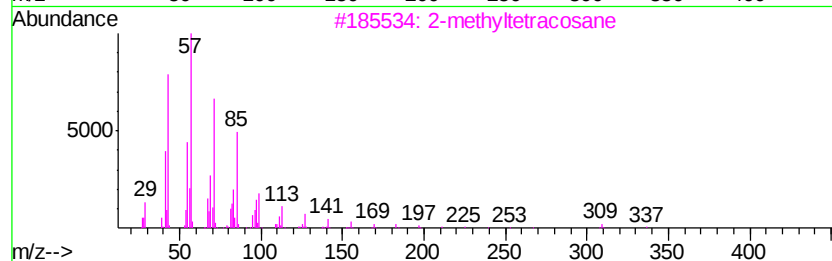
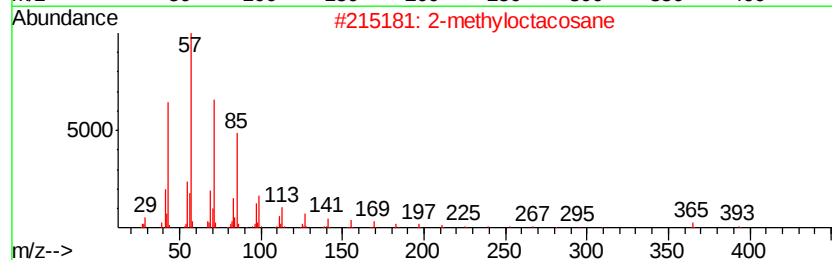
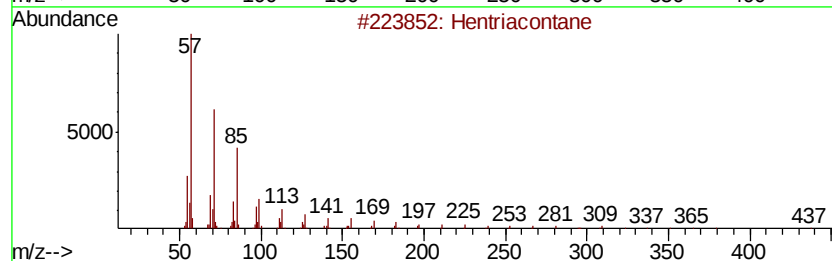
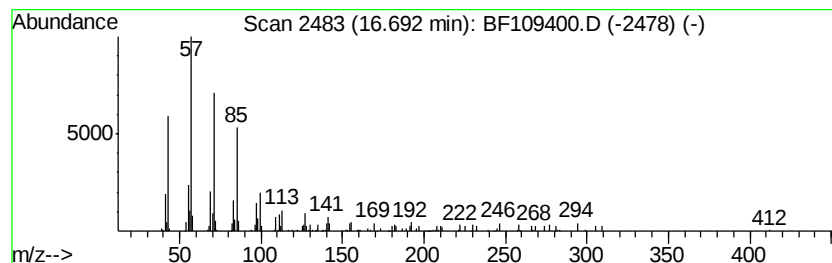
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hentriacontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	7.11 ng	183325	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	87
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		2-methyltetracosane	352	C25H52	1000376-72-6	86
4		Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	86
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
Operator : JU/SJ
Sample : J5077-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PARCEL-193-092418-1

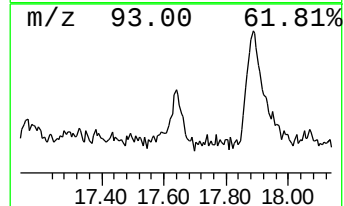
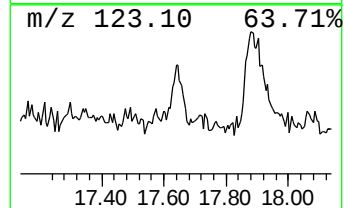
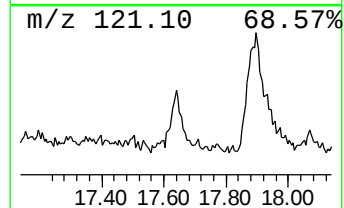
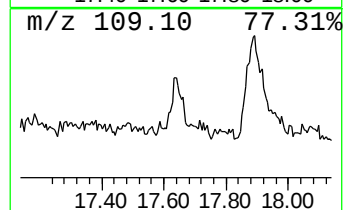
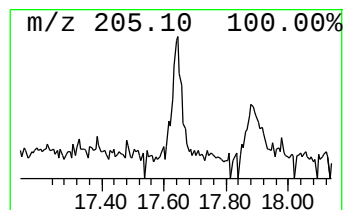
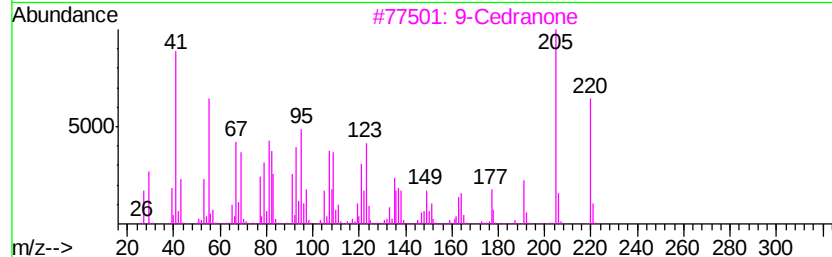
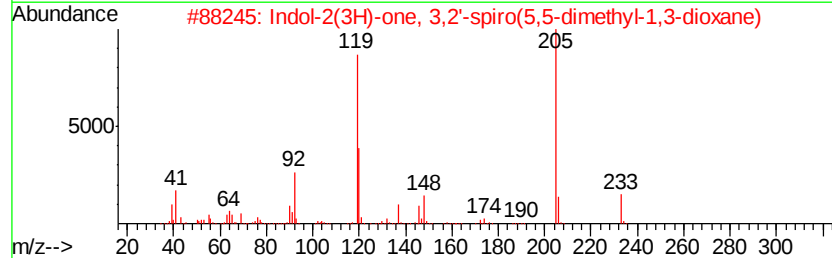
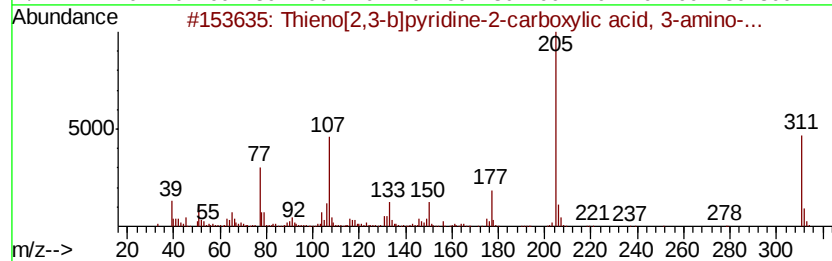
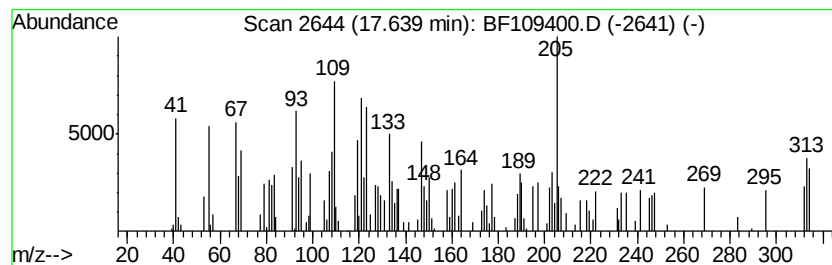
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.64 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.64	3.78 ng	97476	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Thieno[2,3-b]pyridine-2-carboxyl...	311	C17H17N3OS	1000302-46-4	27
2		Indol-2(3H)-one, 3,2'-spiro(5,5-...	233	C13H15N03	1000272-73-5	22
3		9-Cedranone	220	C15H24O	1000156-23-2	20
4		3-(3-Imino-1-pyrazolidinyl)benzo...	205	C10H11N3O2	083671-99-2	20
5		7-Chloro-2,3-dihydrofuro(2,3-b)q...	205	C11H8ClNO	064124-91-0	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
Data File : BF109400.D
Acq On : 27 Sep 2018 21:02
Operator : JU/SJ
Sample : J5077-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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PARCEL-193-092418-1

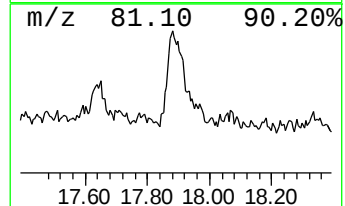
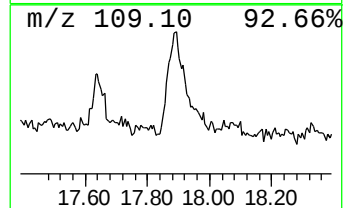
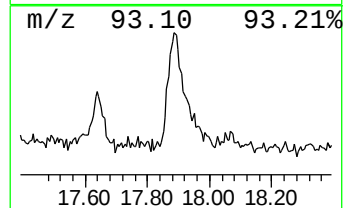
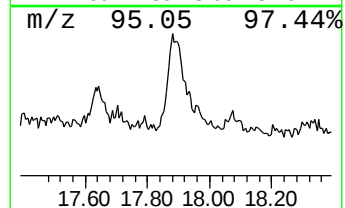
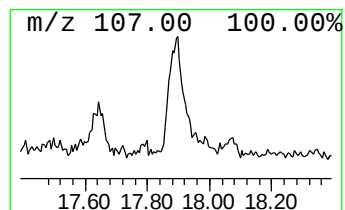
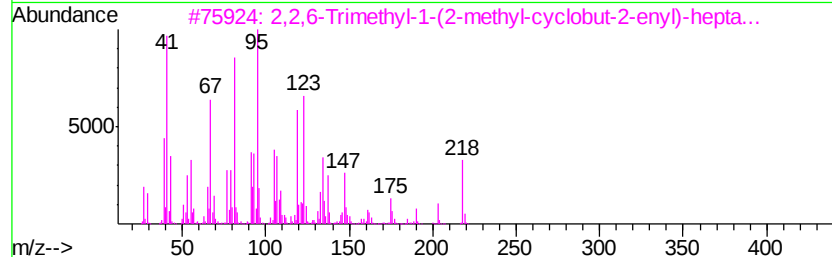
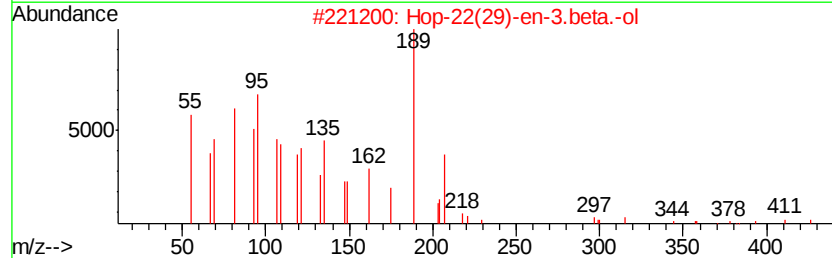
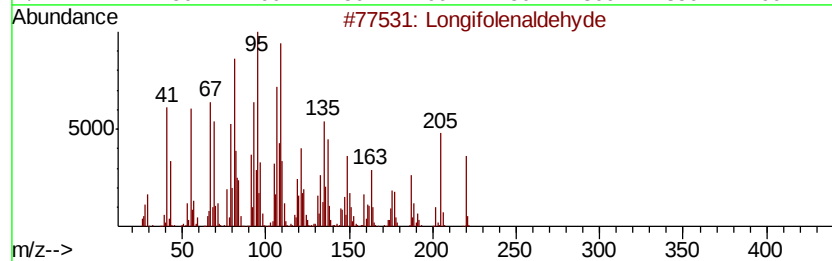
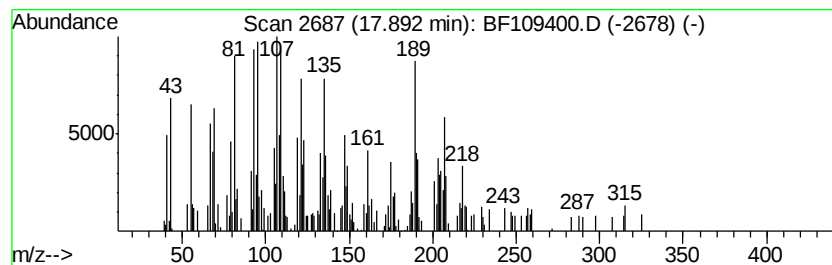
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Longifolenaldehyde Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.89	21.34 ng	550018	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Longifolenaldehyde	220	C15H24O	019890-84-7	59
2		Hop-22(29)-en-3.beta.-ol	426	C30H50O	058801-23-3	55
3		2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	53
4		1,5-Cycloundecadiene, 9-(1-methy...	190	C14H22	062338-55-0	49
5		1,3-Cyclopentadiene, 2,3,4,5-tet...	190	C14H22	1000161-28-7	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092718\
 Data File : BF109400.D
 Acq On : 27 Sep 2018 21:02
 Operator : JU/SJ
 Sample : J5077-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.89	2.3	ng	75179	1	6.70	658301	20.0
unknown6.47	6.47	53.9	ng	1772410	1	6.70	658301	20.0
n-Hexadecanoic acid	11.75	3.9	ng	148955	4	11.22	755293	20.0
Eicosyl pentafluo...	13.71	5.9	ng	257068	5	13.84	874155	20.0
Cyclopentadecane	14.34	9.2	ng	403734	5	13.84	874155	20.0
Supraene	14.69	5.7	ng	147908	6	15.24	515528	20.0
1,19-Eicosadiene	14.75	16.8	ng	431970	6	15.24	515528	20.0
Dodecane, 2-methyl-	14.93	9.2	ng	237952	6	15.24	515528	20.0
Behenic alcohol	14.97	17.9	ng	460576	6	15.24	515528	20.0
Benzo[e]pyrene	15.12	2.2	ng	56490	6	15.24	515528	20.0
Octadecanal	15.47	7.8	ng	201990	6	15.24	515528	20.0
Heptadecane	15.69	14.0	ng	360030	6	15.24	515528	20.0
Heptadecanal	16.42	5.0	ng	128957	6	15.24	515528	20.0
Hentriacontane	16.69	7.1	ng	183325	6	15.24	515528	20.0
unknown17.64	17.64	3.8	ng	97476	6	15.24	515528	20.0
Longifolenaldehyde	17.89	21.3	ng	550018	6	15.24	515528	20.0