

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110718\
 Data File : BF110567.D
 Acq On : 7 Nov 2018 20:35
 Operator : JU/SJ
 Sample : J5813-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-H

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-BF102318.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	2.263	25	30	52	rVB	240179	382077	22.89%	2.363%
2	5.186	524	527	542	rBV	45416	67780	4.06%	0.419%
3	5.575	589	593	607	rBV	1624224	1526510	91.46%	9.439%
4	6.575	759	763	784	rBV	1426021	1669053	100.00%	10.321%
5	6.722	784	788	791	rBV	1916774	1608855	96.39%	9.948%
6	6.951	824	827	833	rBV	532086	440389	26.39%	2.723%
7	7.110	849	854	867	rBV	1311544	1242034	74.42%	7.680%
8	7.516	918	923	941	rBV	1013400	985173	59.03%	6.092%
9	8.233	1041	1045	1062	rBV	486496	518340	31.06%	3.205%
10	9.304	1223	1227	1238	rBV	1470230	1487249	89.11%	9.196%
11	9.698	1291	1294	1306	rVB	167227	156976	9.41%	0.971%
12	9.992	1340	1344	1354	rBV	501942	479905	28.75%	2.968%
13	10.786	1475	1479	1486	rBV	744682	706109	42.31%	4.366%
14	11.486	1595	1598	1601	rBV	423008	405117	24.27%	2.505%
15	11.510	1601	1602	1609	rVB	98739	87618	5.25%	0.542%
16	11.992	1681	1684	1688	rBV	89424	89981	5.39%	0.556%
17	12.704	1802	1805	1811	rBV	199920	186740	11.19%	1.155%
18	12.915	1838	1841	1842	rBV	28312	22119	1.33%	0.137%
19	12.939	1842	1845	1850	rVB	153949	141430	8.47%	0.875%
20	13.068	1863	1867	1870	rBV	1387334	1159069	69.44%	7.167%
21	13.262	1893	1900	1904	rBV4	25087	42964	2.57%	0.266%
22	13.533	1943	1946	1951	rVB	113515	97377	5.83%	0.602%
23	13.886	2003	2006	2009	rBV2	20818	25404	1.52%	0.157%
24	13.945	2013	2016	2021	rVV2	151101	192361	11.53%	1.189%
25	13.986	2021	2023	2030	rVB	28618	37391	2.24%	0.231%
26	14.139	2044	2049	2052	rBV	535859	589408	35.31%	3.645%
27	14.162	2052	2053	2059	rVB	89354	69627	4.17%	0.431%
28	14.262	2065	2070	2073	rBV4	31766	48285	2.89%	0.299%
29	14.568	2119	2122	2130	rBV5	50229	85457	5.12%	0.528%
30	14.886	2173	2176	2180	rVB	57195	59838	3.59%	0.370%
31	15.039	2199	2202	2207	rVB	293205	288731	17.30%	1.785%
32	15.209	2228	2231	2233	rBV	61767	82785	4.96%	0.512%
33	15.239	2233	2236	2238	rVV	129719	155537	9.32%	0.962%
34	15.268	2238	2241	2248	rVB	162904	221447	13.27%	1.369%

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

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

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35	15.603	2295	2298	2300	rBV	39037	52335	3.14%	0.324%
36	15.633	2300	2303	2305	rVV	81137	98957	5.93%	0.612%
37	15.668	2305	2309	2317	rVB	233160	325864	19.52%	2.015%
38	15.851	2337	2340	2347	rVB3	38977	60956	3.65%	0.377%
39	16.086	2376	2380	2386	rVB	122136	199557	11.96%	1.234%
40	16.621	2467	2471	2480	rVB	45747	75203	4.51%	0.465%

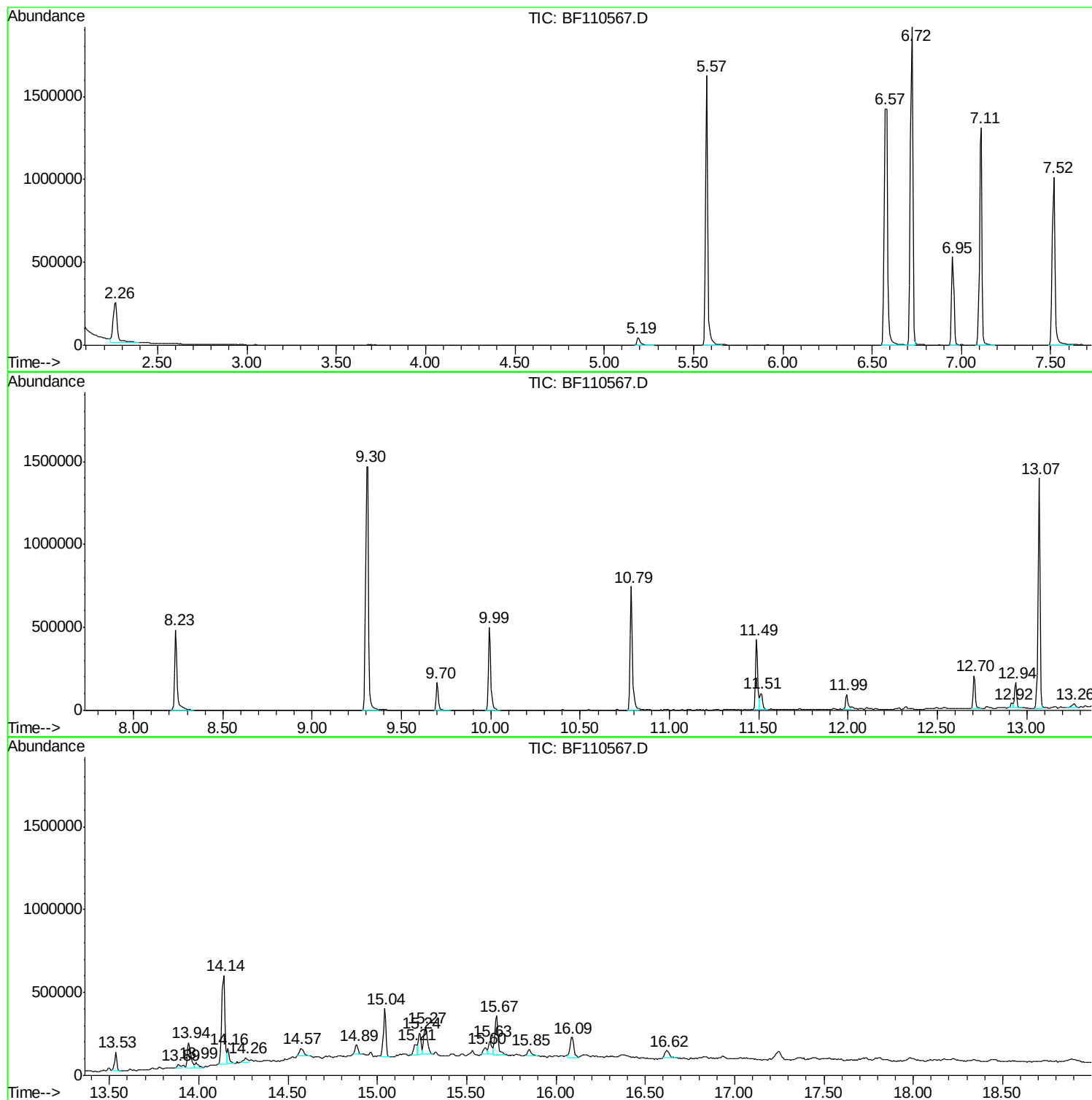
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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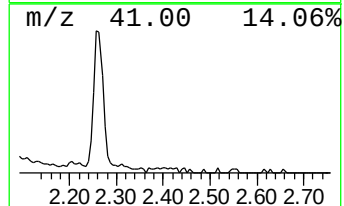
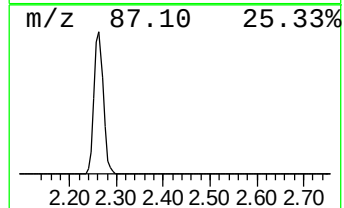
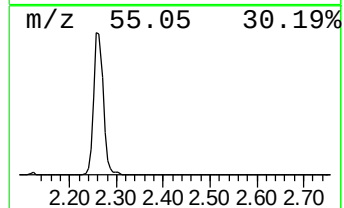
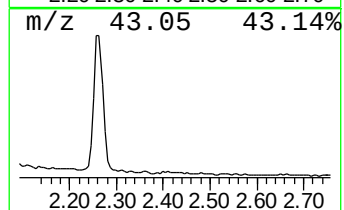
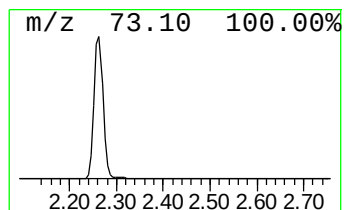
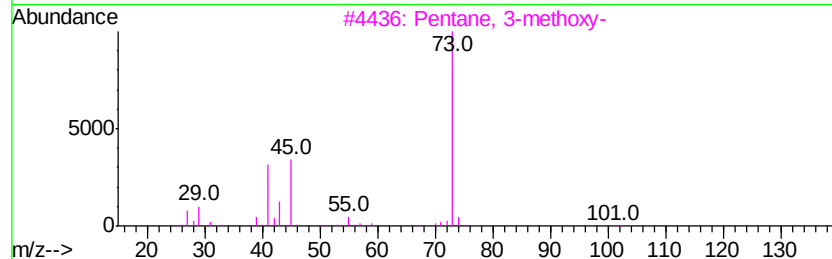
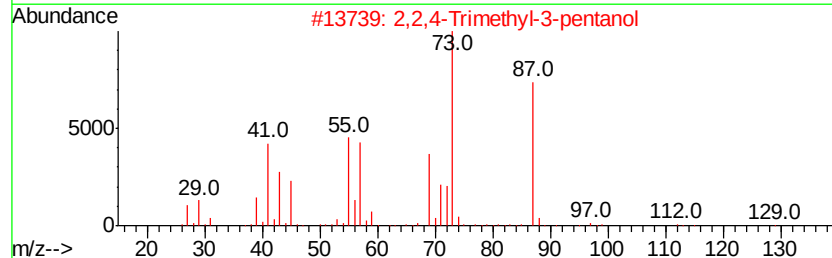
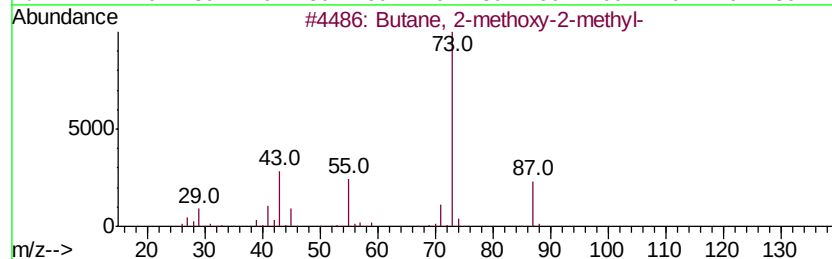
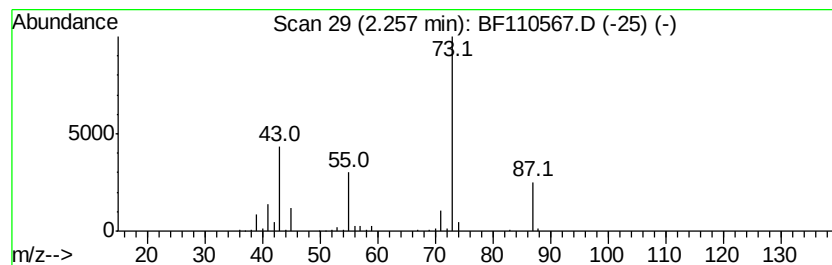
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.26	17.35 ng	382077	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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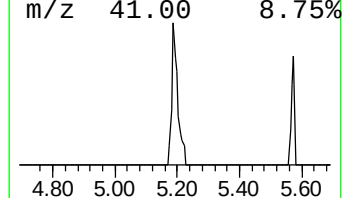
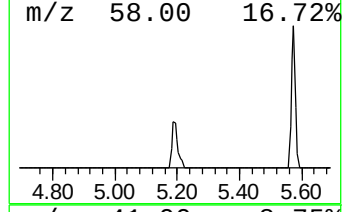
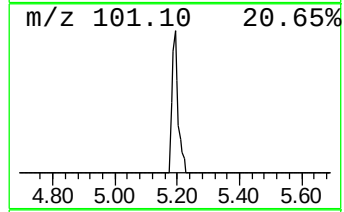
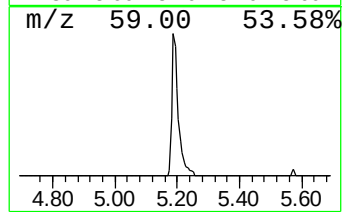
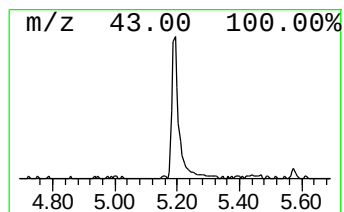
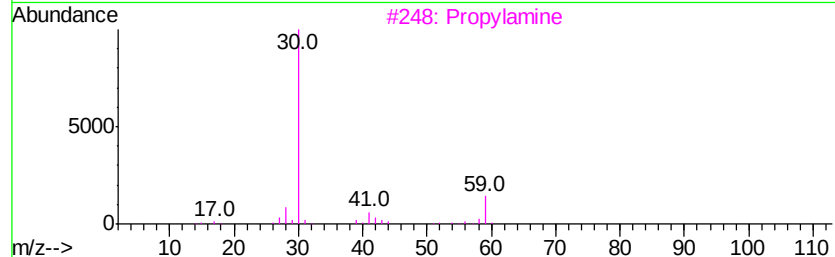
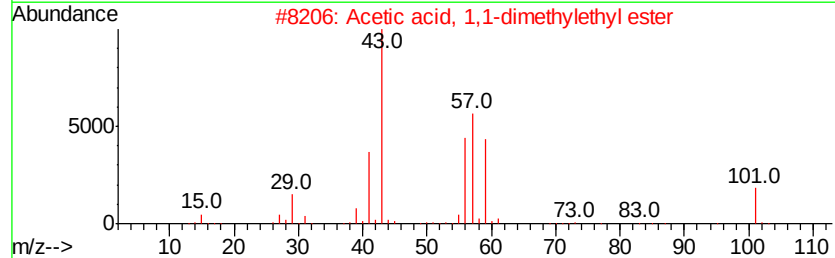
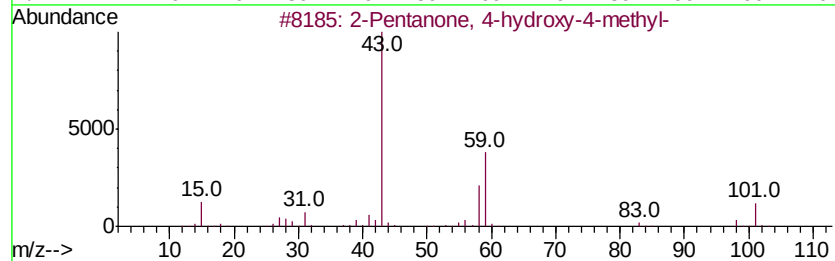
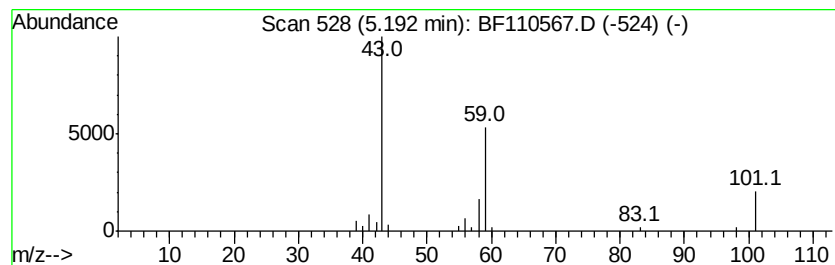
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TIC Library : C:\DATABASE\NIST11.L
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 Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.08 ng	67780	1,4-Dichlorobenzene-d4	6.95

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			Propylamine	59	C3H9N	000107-10-8	9
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
5			3-Hexanol	102	C6H14O	000623-37-0	9



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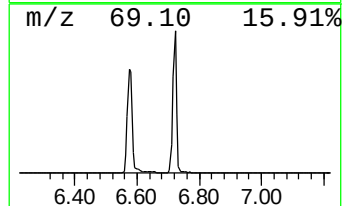
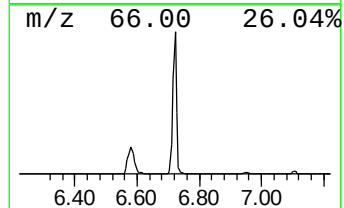
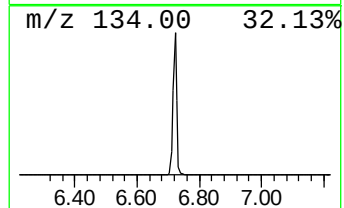
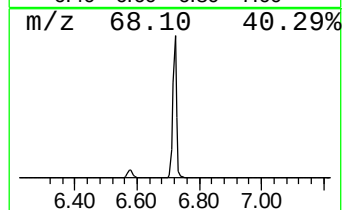
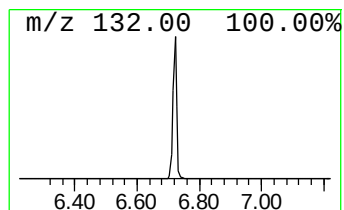
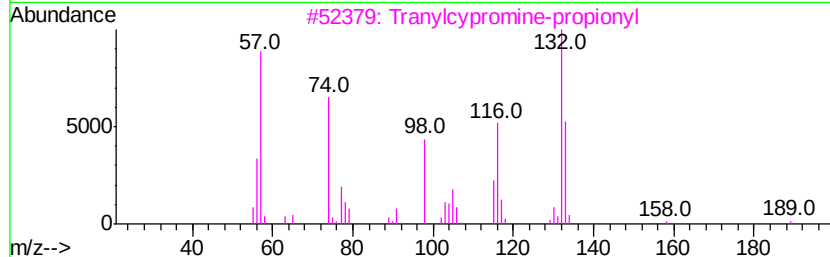
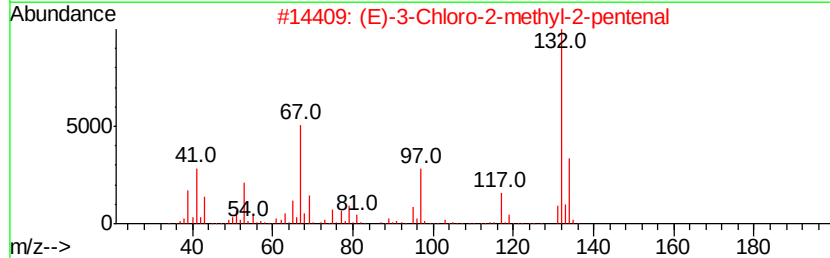
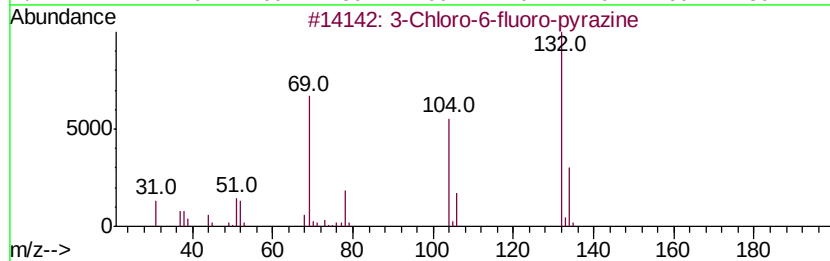
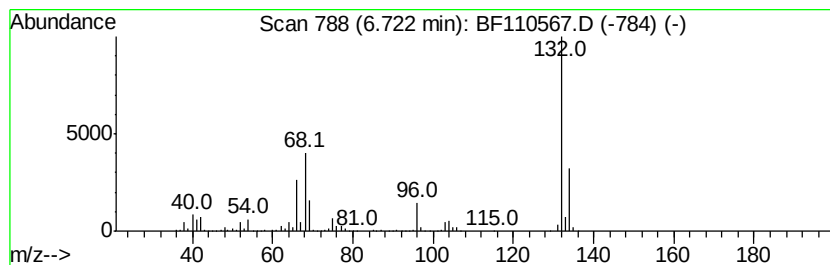
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.72 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	73.07 ng	1608860	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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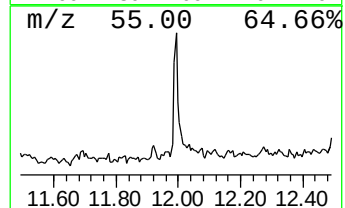
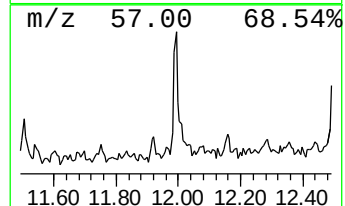
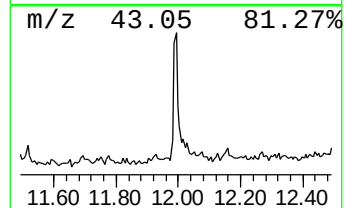
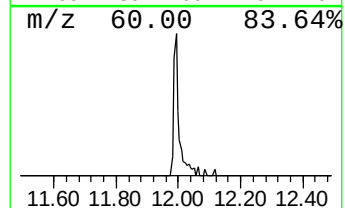
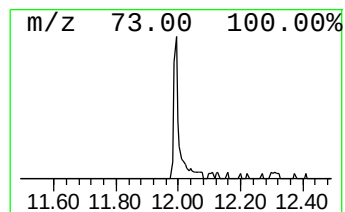
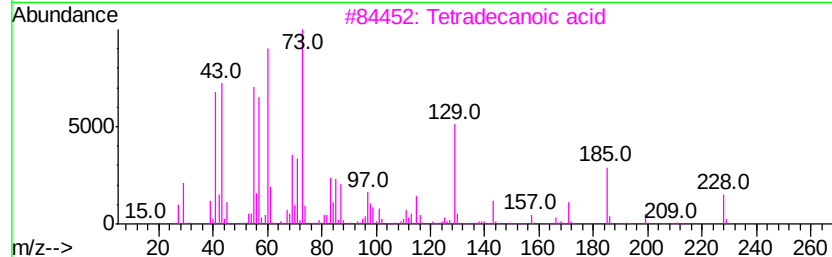
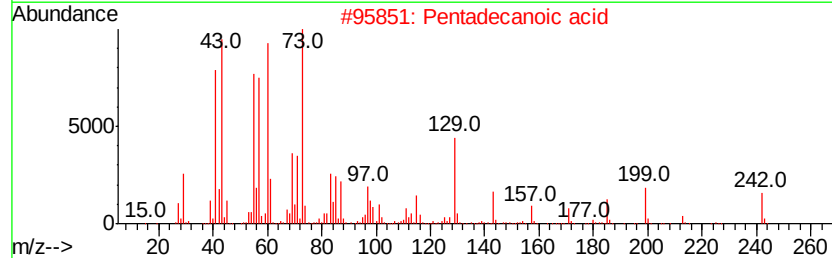
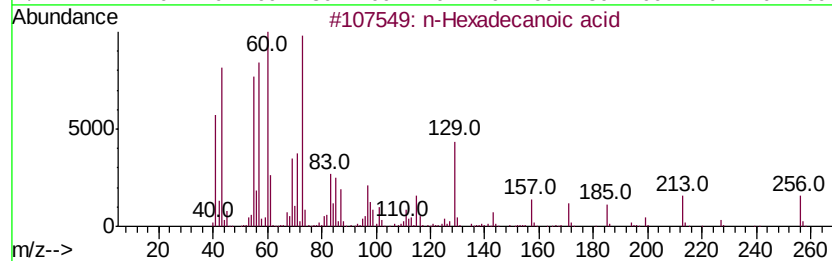
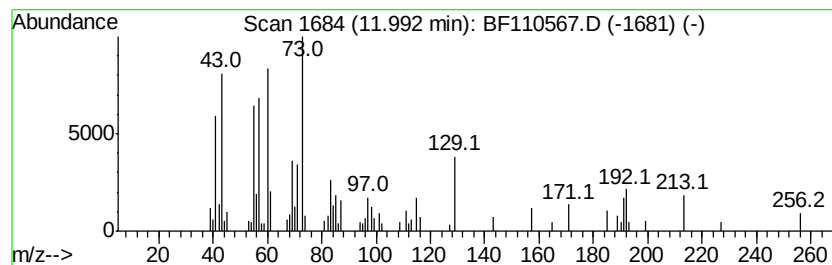
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.99	4.44 ng	89981	Phenanthrene-d10		11.49	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	70
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4		Tridecanoic acid	214	C13H26O2	000638-53-9	62
5		n-Decanoic acid	172	C10H20O2	000334-48-5	53



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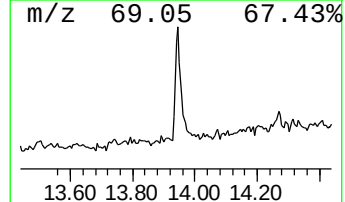
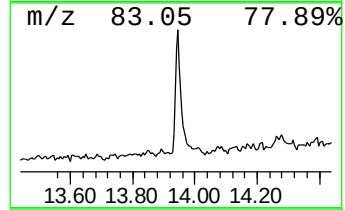
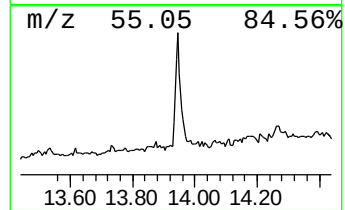
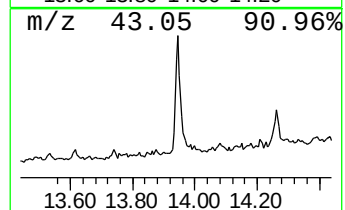
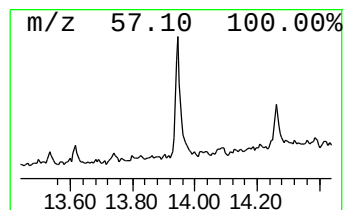
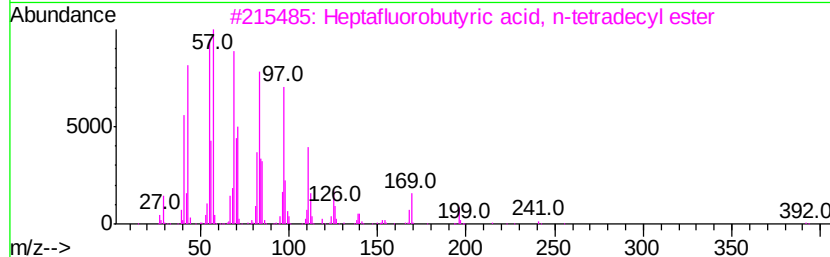
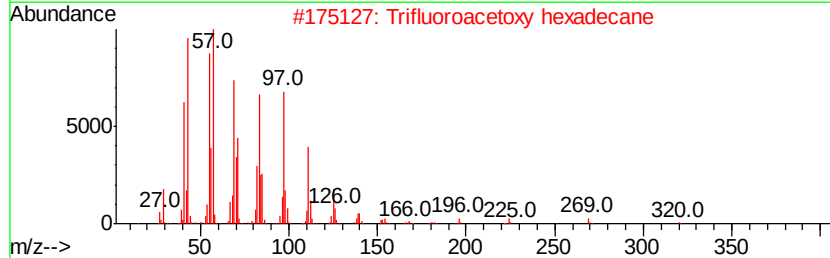
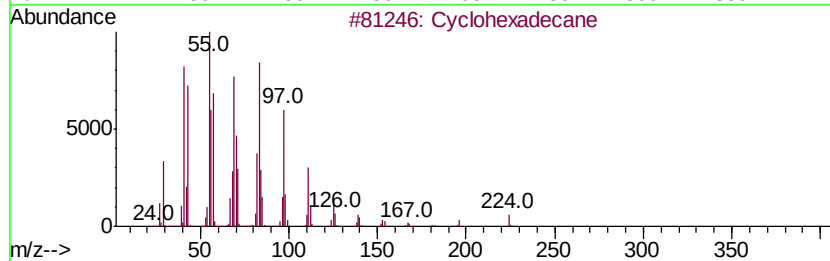
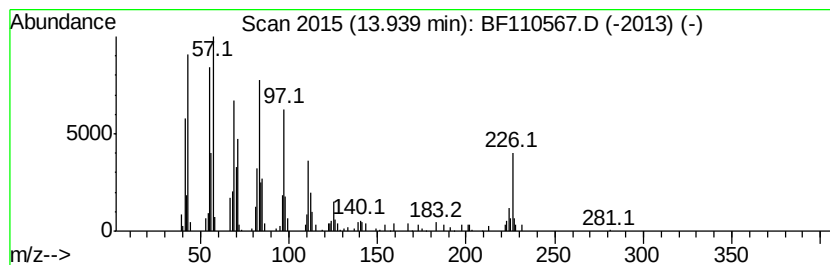
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclohexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	6.53 ng	192361	Chrysene-d12	14.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 98
2		Trifluoroacetoxy hexadecane	338 C18H33F3O2	006222-03-3 93
3		Heptafluorobutyric acid, n-tetra...	410 C18H29F7O2	007365-36-8 93
4		Heptadecyl heptafluorobutyrate	452 C21H35F7O2	959085-66-6 93
5		Octadecyl trifluoroacetate	366 C20H37F3O2	079392-43-1 90



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Acq On : 7 Nov 2018 20:35
Operator : JU/SJ
Sample : J5813-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

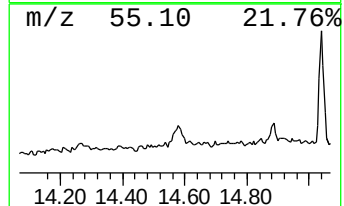
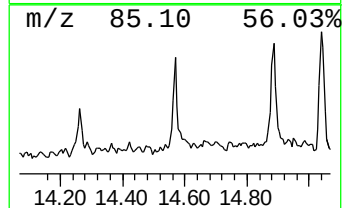
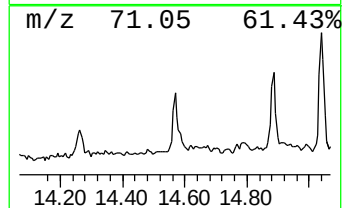
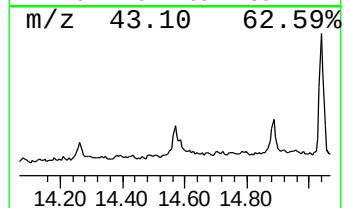
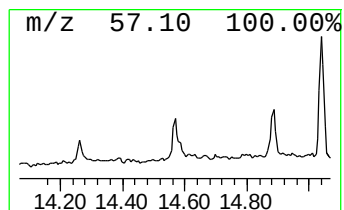
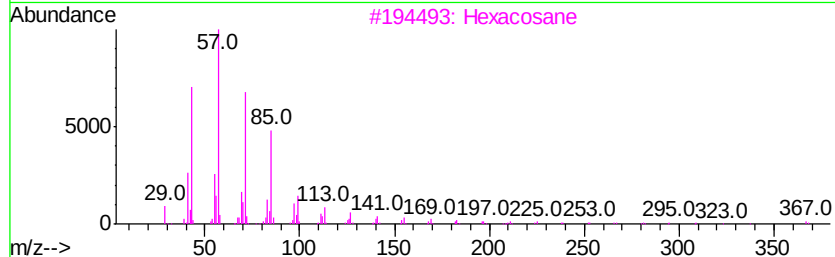
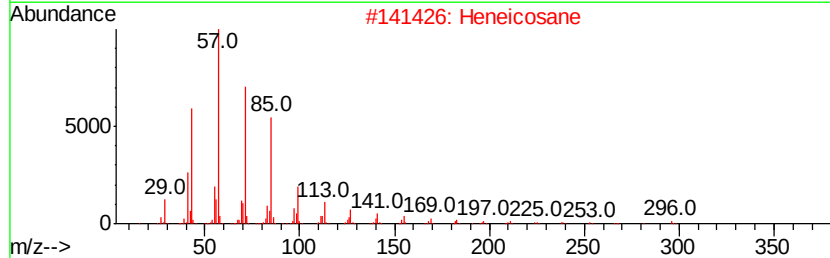
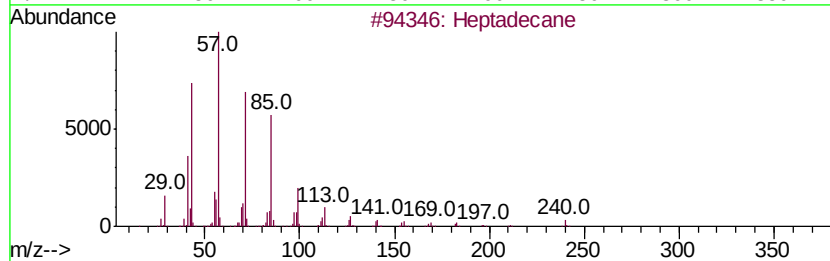
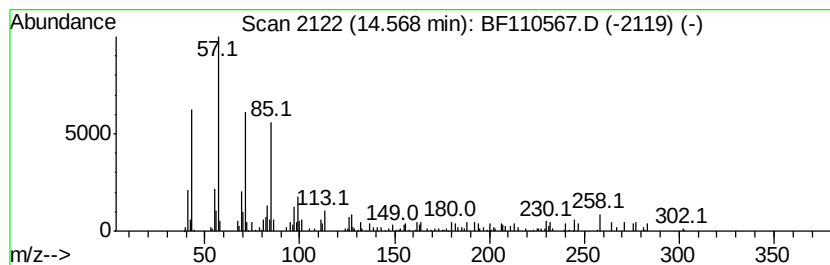
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.57	2.90 ng	85457	Chrysene-d12	14.14	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	86
2	Heneicosane	296	C21H44	000629-94-7	64
3	Hexacosane	366	C26H54	000630-01-3	64
4	Tricosane, 2-methyl-	338	C24H50	001928-30-9	52
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110718\
Data File : BF110567.D
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Operator : JU/SJ
Sample : J5813-08
Misc :
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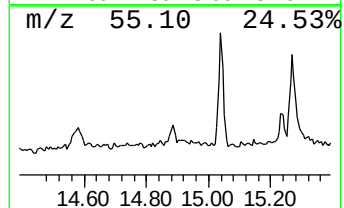
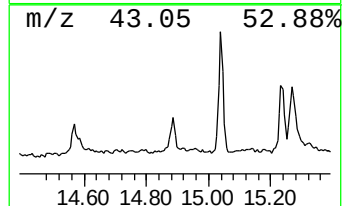
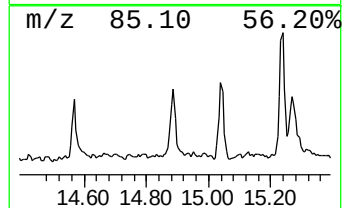
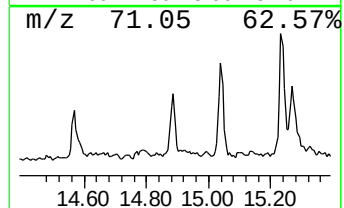
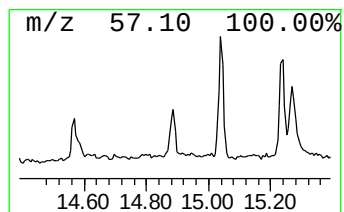
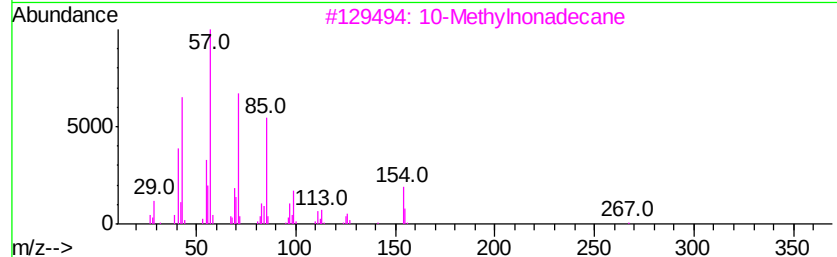
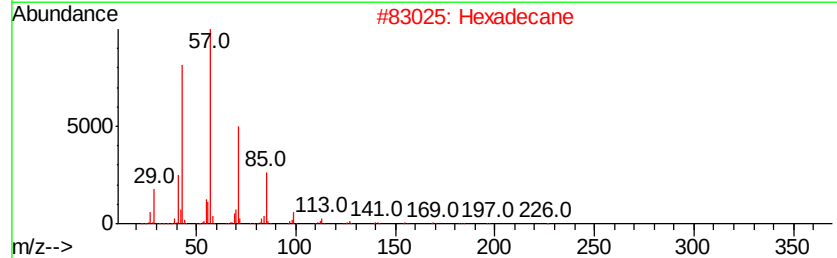
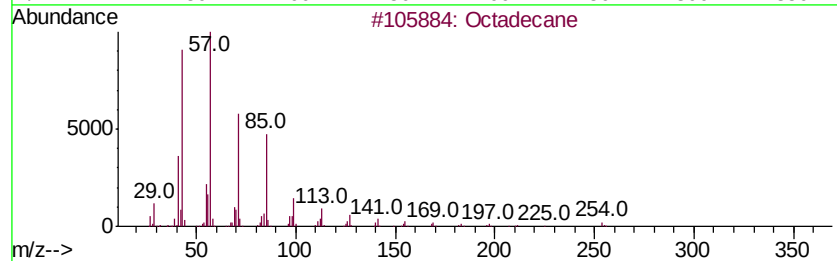
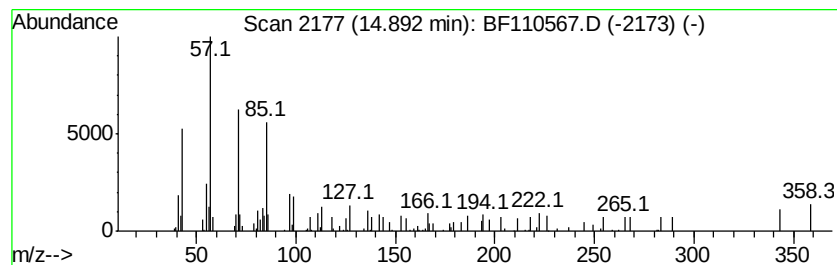
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	2.03 ng	59838	Chrysene-d12	14.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octadecane	254 C18H38	000593-45-3	83
2	Hexadecane	226 C16H34	000544-76-3	78
3	10-Methylnonadecane	282 C20H42	056862-62-5	53
4	Heptacosane	380 C27H56	000593-49-7	52
5	Octacosane	394 C28H58	000630-02-4	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110718\
Data File : BF110567.D
Acq On : 7 Nov 2018 20:35
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Sample : J5813-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

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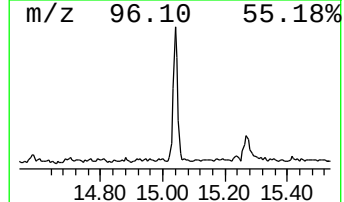
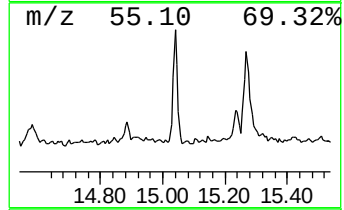
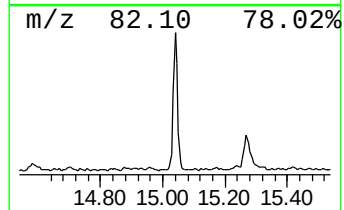
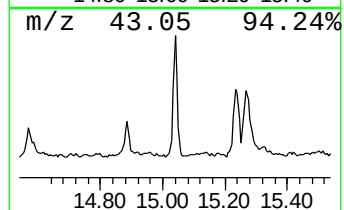
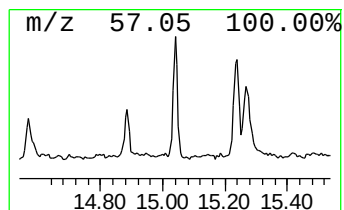
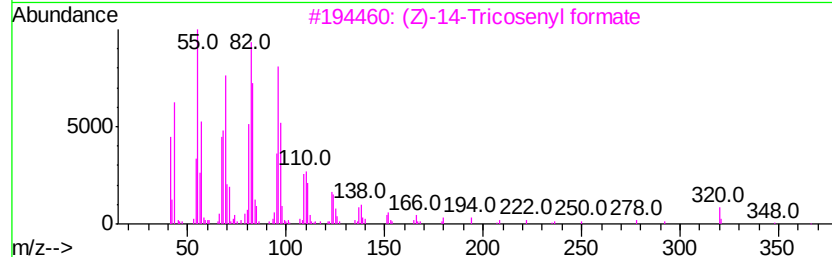
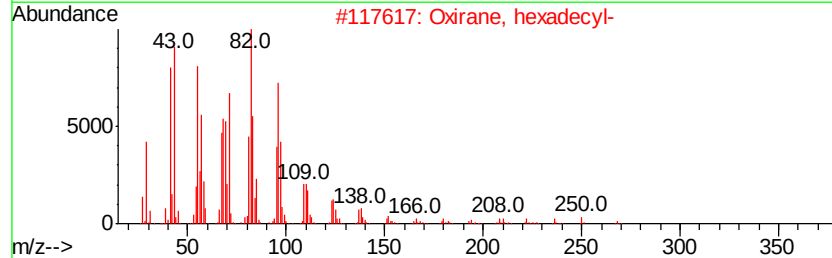
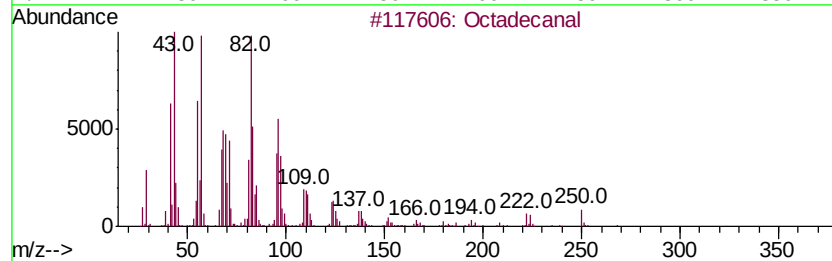
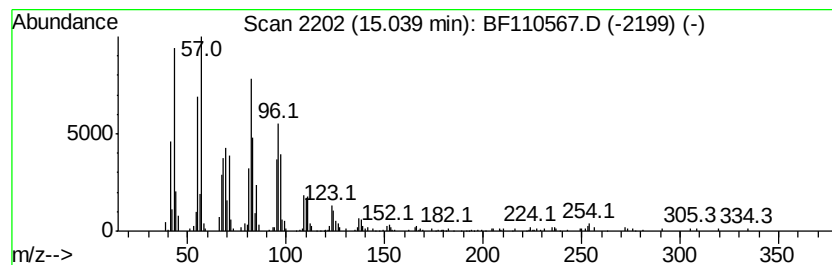
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanal Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.04	17.72 ng	288731	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanal	268	C18H36O	000638-66-4	91
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3		(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	90
4		16-Heptadecenal	252	C17H32O	1000144-57-9	87
5		1,19-Eicosadiene	278	C20H38	014811-95-1	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110718\
Data File : BF110567.D
Acq On : 7 Nov 2018 20:35
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Sample : J5813-08
Misc :
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Instrument :
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ClientSampleId :
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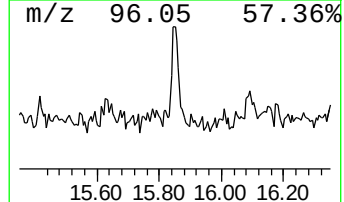
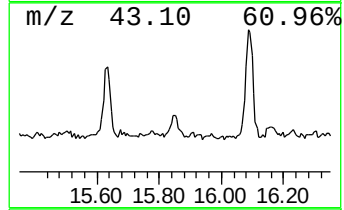
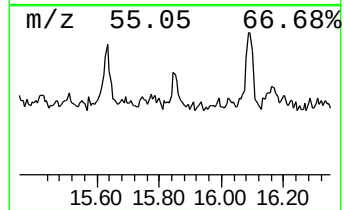
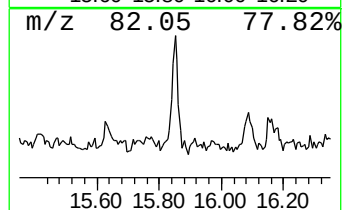
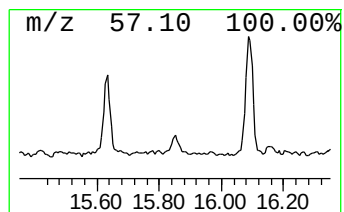
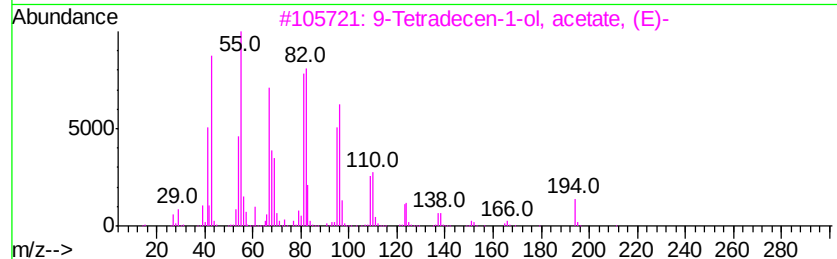
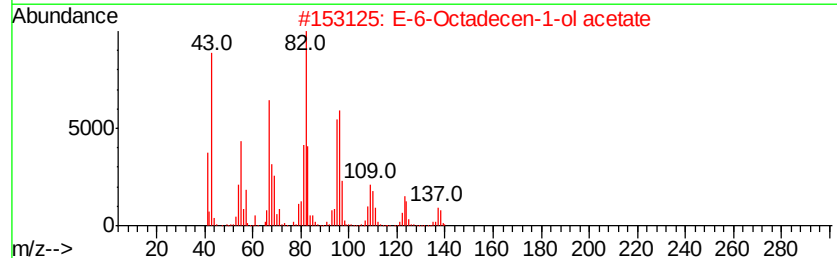
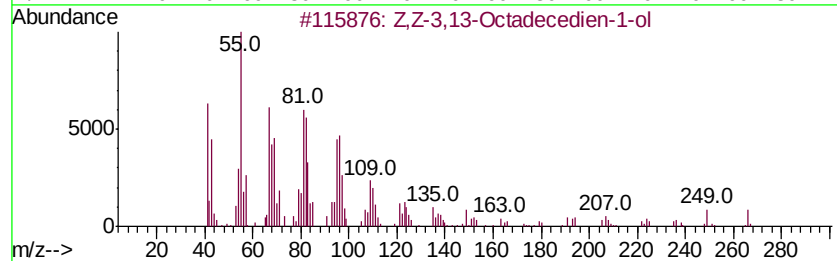
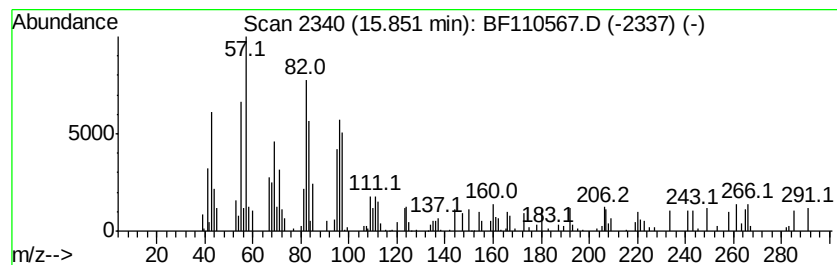
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Z,Z-3,13-Octadecadien-1-ol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.85	3.74 ng	60956	Perylene-d12	15.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z,Z-3,13-Octadecadien-1-ol	266	C18H34O	1000131-10-7	64
2			E-6-Octadecen-1-ol acetate	310	C20H38O2	1000130-94-7	50
3			9-Tetradecen-1-ol, acetate, (E)-	254	C16H30O2	023192-82-7	47
4			1-Oxacyclopentadecan-2-one, 15-i...	266	C17H30O2	089328-44-9	46
5			Cyclohexane, 1,1'-methylenebis-	180	C13H24	003178-23-2	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110718\
Data File : BF110567.D
Acq On : 7 Nov 2018 20:35
Operator : JU/SJ
Sample : J5813-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

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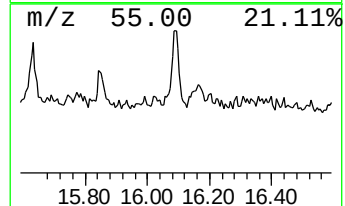
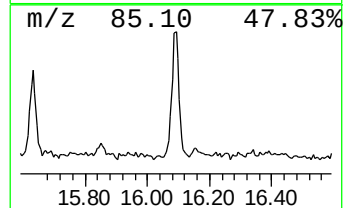
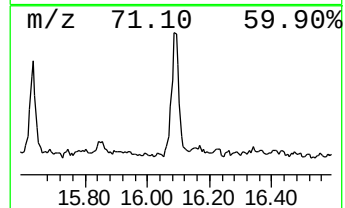
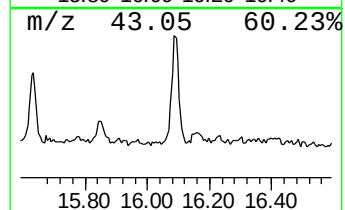
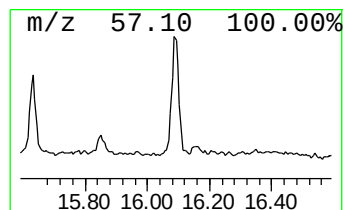
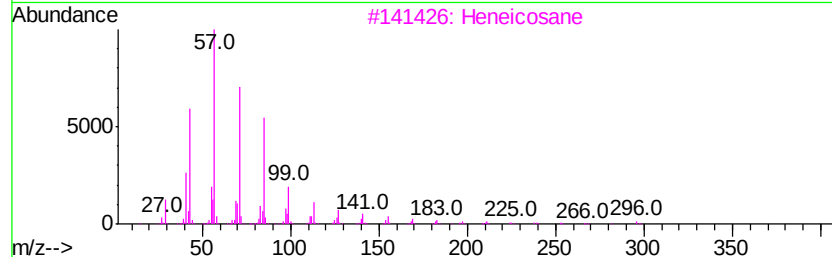
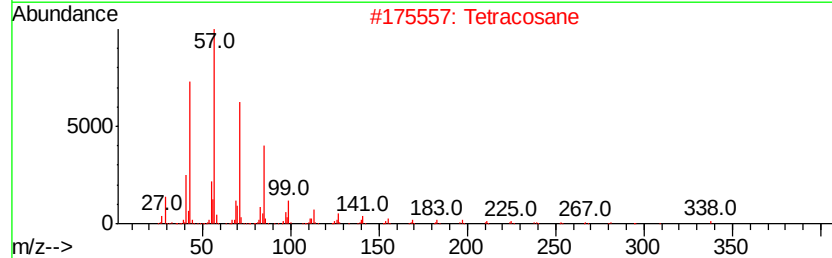
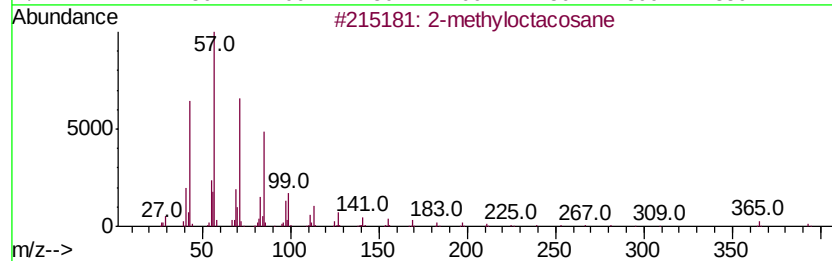
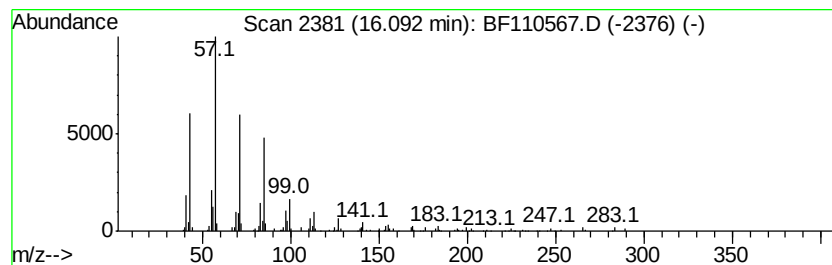
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-methyloctacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	12.25 ng	199557	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	90
2		Tetracosane	338	C24H50	000646-31-1	87
3		Heneicosane	296	C21H44	000629-94-7	74
4		Hexadecane	226	C16H34	000544-76-3	72
5		Octadecane	254	C18H38	000593-45-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110718\
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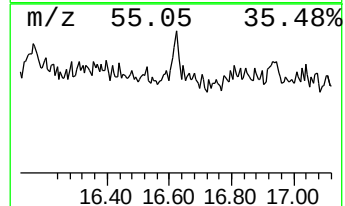
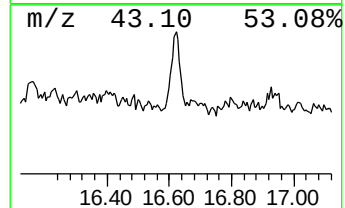
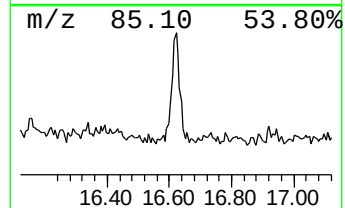
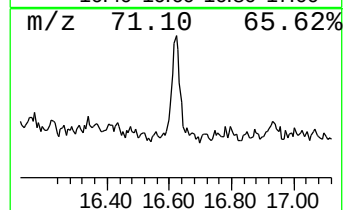
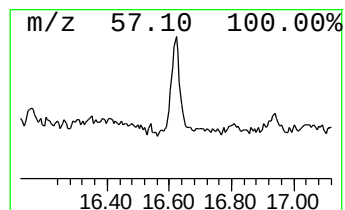
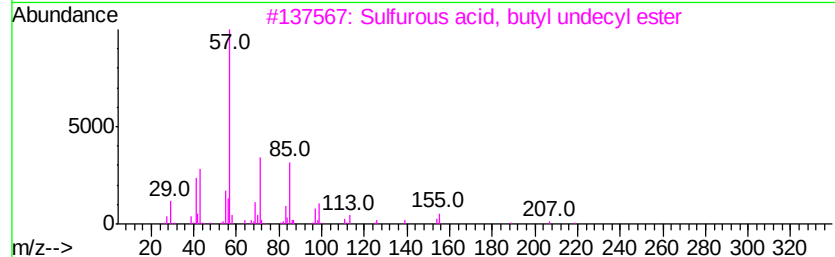
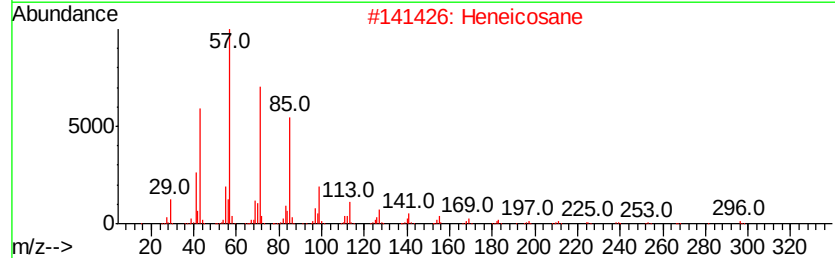
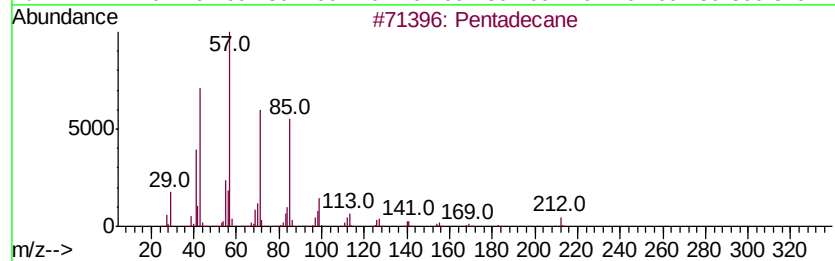
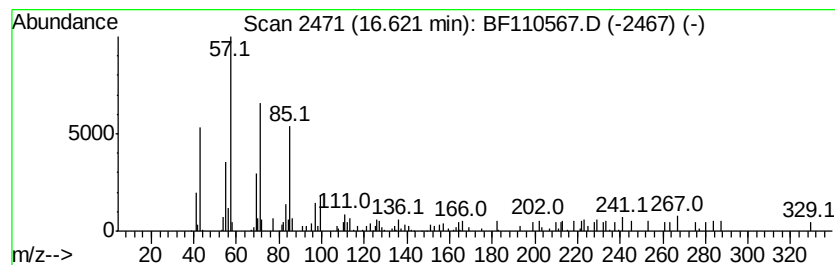
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.62	4.62 ng	75203	Perylene-d12	15.67
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentadecane	212 C15H32	000629-62-9	58
2	Heneicosane	296 C21H44	000629-94-7	52
3	Sulfurous acid, butyl undecyl ester	292 C15H32O3S	1000309-17-8	50
4	Sulfurous acid, 2-ethylhexyl hex...	278 C14H30O3S	1000309-20-2	38
5	Tridecane, 2-methyl-	198 C14H30	001560-96-9	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110718\
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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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
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2-Pentanone, 4-hy...	5.19	3.1	ng	67780	1	6.95	440389	20.0
unknown6.72	6.72	73.1	ng	1608860	1	6.95	440389	20.0
n-Hexadecanoic acid	11.99	4.4	ng	89981	4	11.49	405117	20.0
Cyclohexadecane	13.94	6.5	ng	192361	5	14.14	589408	20.0
Heptadecane	14.57	2.9	ng	85457	5	14.14	589408	20.0
Octadecane	14.89	2.0	ng	59838	5	14.14	589408	20.0
Octadecanal	15.04	17.7	ng	288731	6	15.67	325864	20.0
Z,Z-3,13-Octadece...	15.85	3.7	ng	60956	6	15.67	325864	20.0
2-methyloctacosane	16.09	12.3	ng	199557	6	15.67	325864	20.0
Pentadecane	16.62	4.6	ng	75203	6	15.67	325864	20.0