

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022819\
 Data File : BF112776.D
 Acq On : 28 Feb 2019 22:12
 Operator : JU/SJ
 Sample : K1645-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SR-9-D

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-BF022719.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	4.299	371	376	385	rBV	104635	138793	3.07%	0.328%
2	4.951	476	487	493	rBV	53897	74908	1.65%	0.177%
3	5.357	550	556	559	rBV	4164617	4136449	91.39%	9.781%
4	6.375	724	729	733	rBV	3799397	4362687	96.39%	10.316%
5	6.516	747	753	756	rBV	4140615	4366853	96.48%	10.326%
6	6.745	788	792	795	rBV	1417051	1238215	27.36%	2.928%
7	6.898	814	818	822	rBV	3670321	3409755	75.33%	8.063%
8	7.304	881	887	890	rBV	2628128	2766565	61.12%	6.542%
9	8.022	1005	1009	1013	rBV	1782486	1506237	33.28%	3.562%
10	9.098	1187	1192	1195	rBV	4983983	4526285	100.00%	10.703%
11	9.486	1255	1258	1262	rBV	289844	243739	5.38%	0.576%
12	9.663	1285	1288	1295	rVB2	30027	57610	1.27%	0.136%
13	9.769	1302	1306	1313	rBV	1577368	1531851	33.84%	3.622%
14	10.557	1435	1440	1443	rBV	2521885	2513752	55.54%	5.944%
15	11.245	1554	1557	1561	rBV	1397520	1324669	29.27%	3.132%
16	11.786	1645	1649	1659	rBV	737887	719131	15.89%	1.700%
17	12.569	1778	1782	1790	rBV	248169	284580	6.29%	0.673%
18	12.704	1798	1805	1811	rBV2	138251	193562	4.28%	0.458%
19	12.833	1822	1827	1830	rBV	3816529	3586091	79.23%	8.480%
20	13.733	1976	1980	1987	rBV	767230	787691	17.40%	1.863%
21	13.874	2000	2004	2007	rBV	1567858	1430810	31.61%	3.383%
22	14.063	2032	2036	2039	rBV	88240	92511	2.04%	0.219%
23	14.363	2084	2087	2093	rBV	176689	177012	3.91%	0.419%
24	14.663	2135	2138	2143	rVB	199537	204751	4.52%	0.484%
25	14.733	2147	2150	2152	rVB	56228	54859	1.21%	0.130%
26	14.980	2188	2192	2198	rBV	220141	251812	5.56%	0.595%
27	15.286	2239	2244	2248	rBV	1237958	1453128	32.10%	3.436%
28	15.339	2249	2253	2258	rVB	204082	235436	5.20%	0.557%
29	15.745	2318	2322	2325	rBV	156628	208989	4.62%	0.494%
30	15.786	2325	2329	2338	rVB	181708	297612	6.58%	0.704%
31	16.215	2399	2402	2407	rVB2	77941	114090	2.52%	0.270%

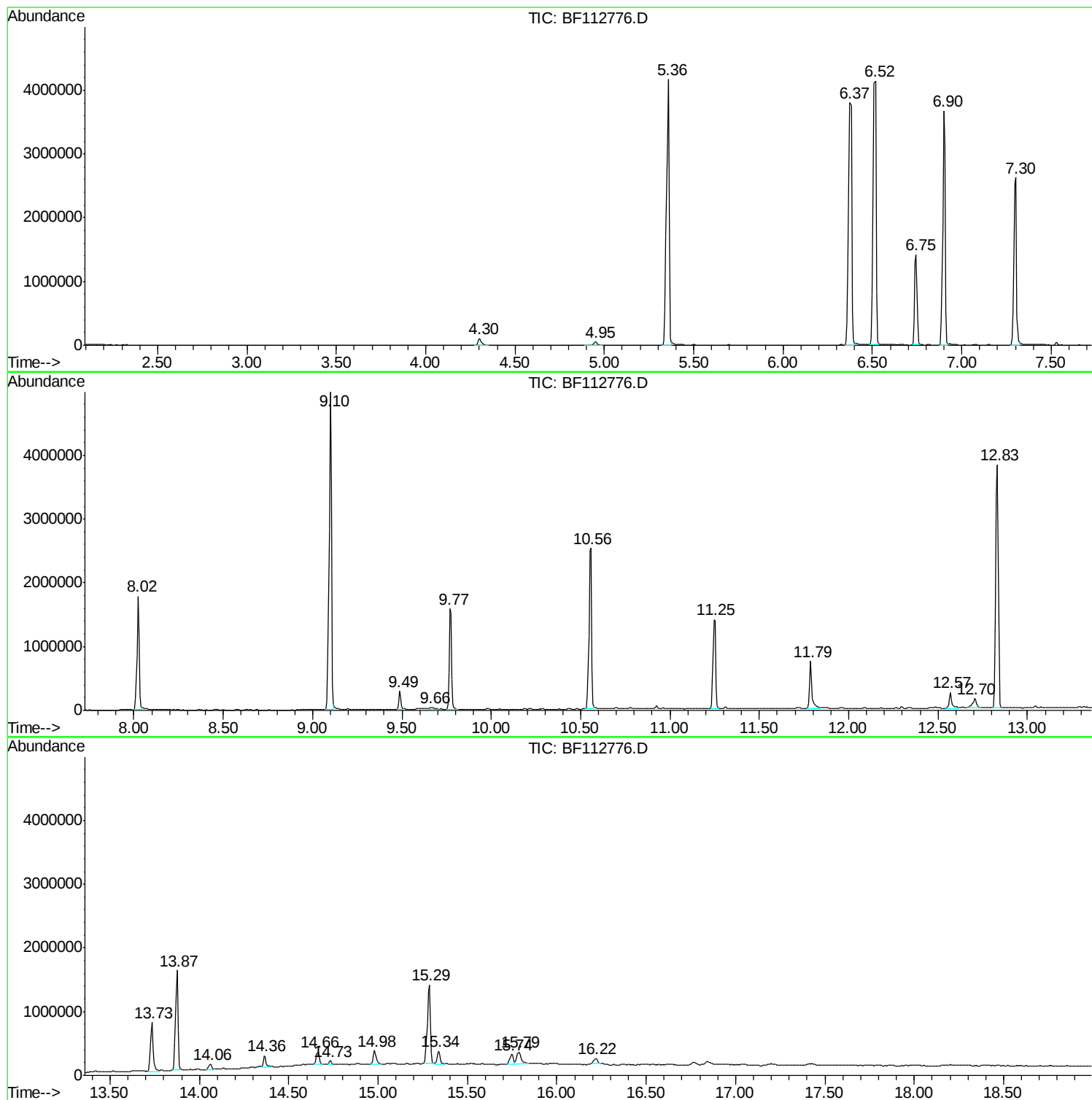
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Instrument :
BNA_F
ClientSampled :
SR-9-D

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.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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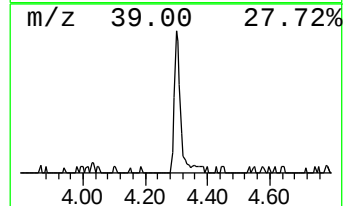
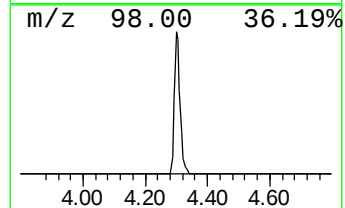
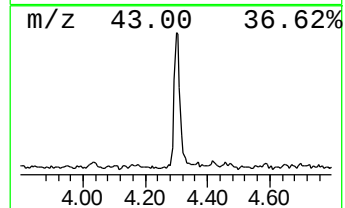
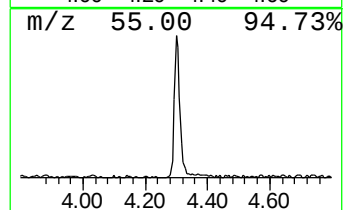
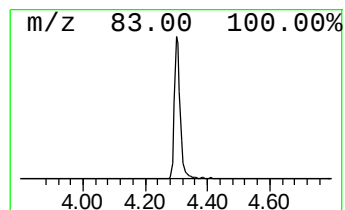
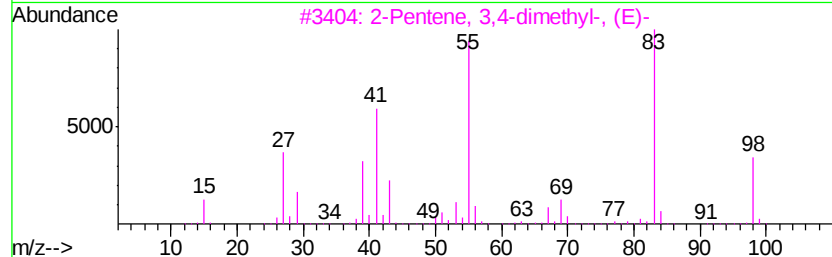
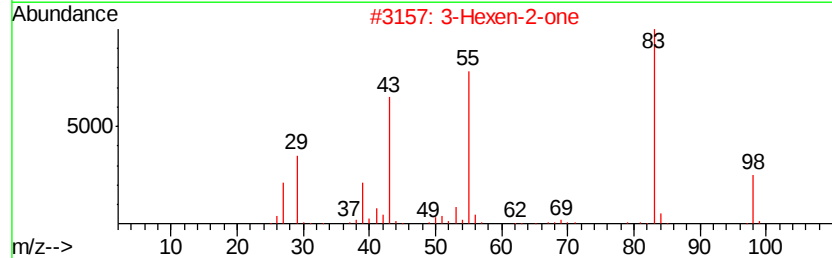
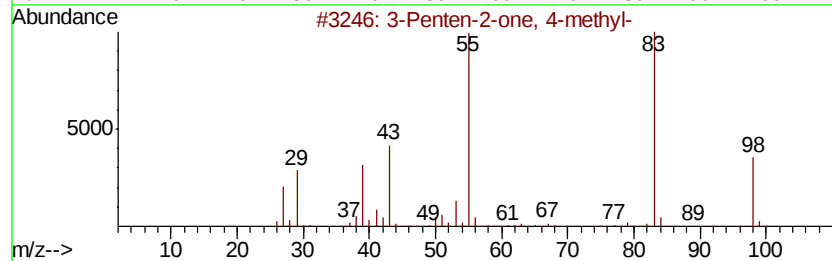
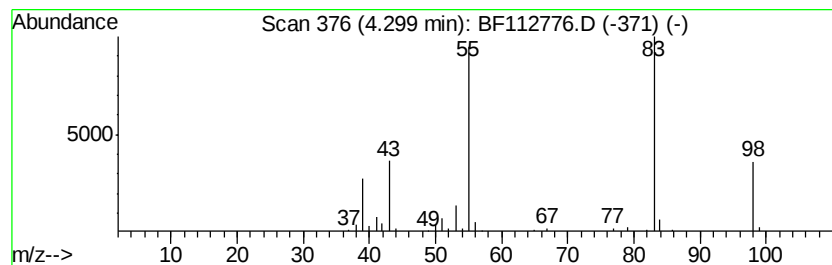
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.30	2.24 ng	138793	1,4-Dichlorobenzene-d4	6.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2			3-Hexen-2-one	98	C6H10O	000763-93-9	91
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	87
5			2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	86



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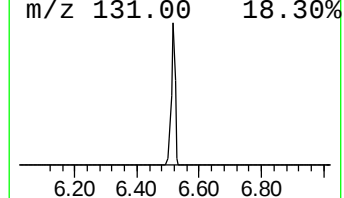
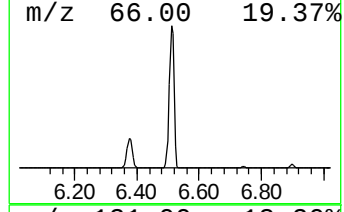
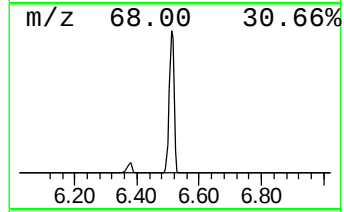
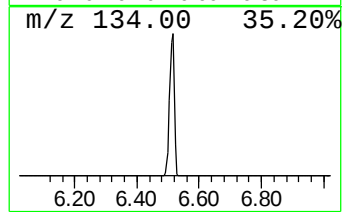
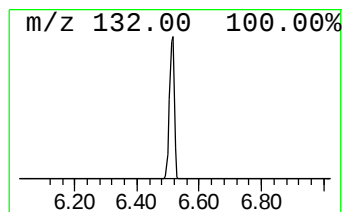
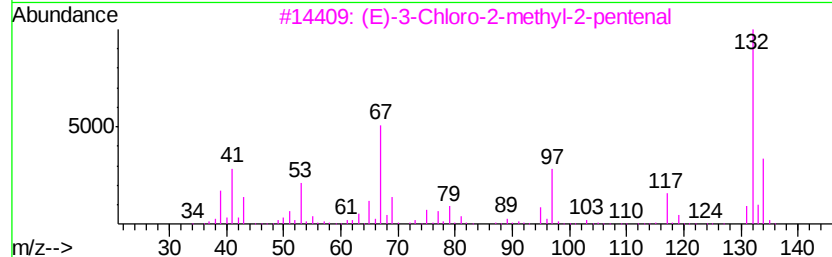
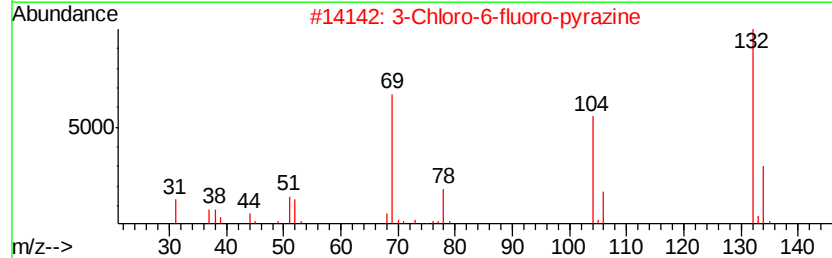
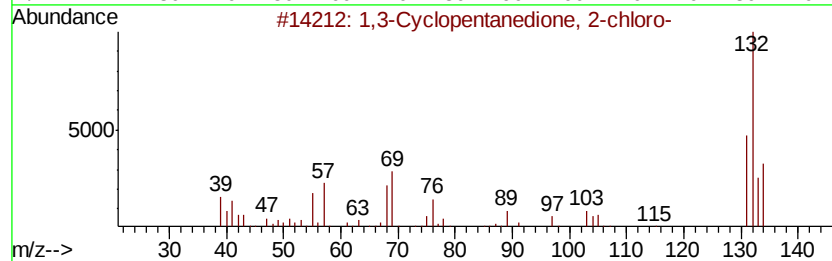
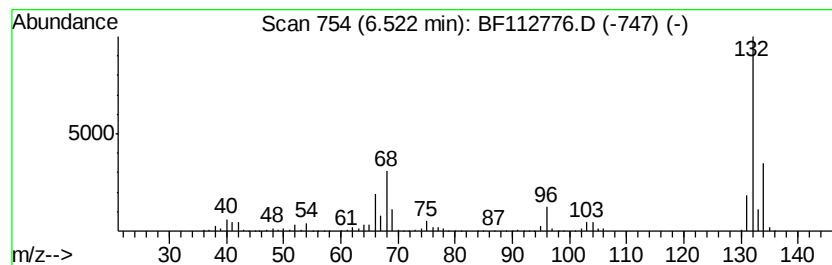
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Peak Number 2 unknown6.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	70.53 ng	4366850	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	47
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Amphetaminil	250	C17H18N2	017590-01-1	10



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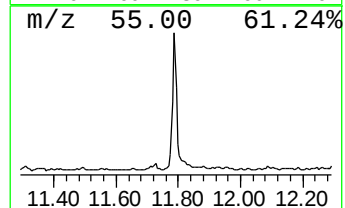
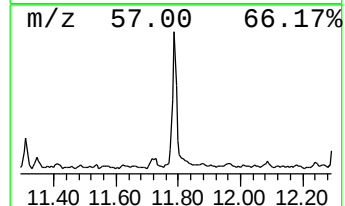
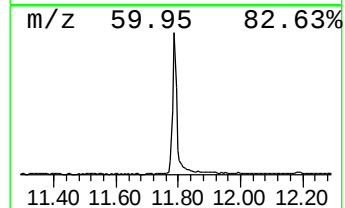
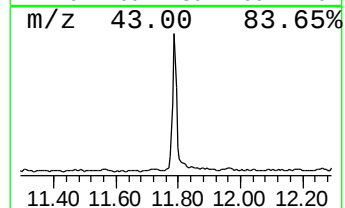
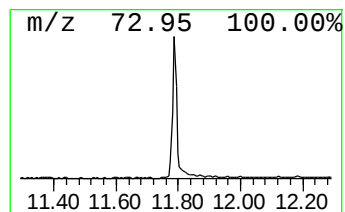
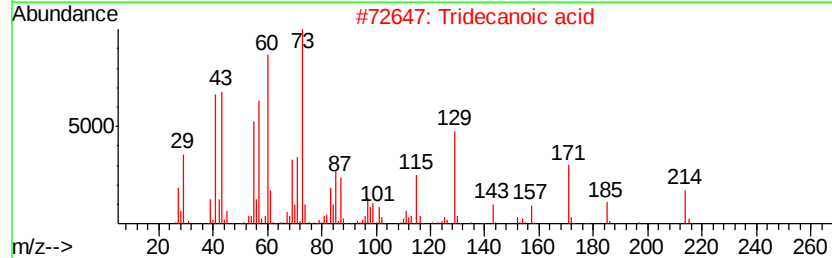
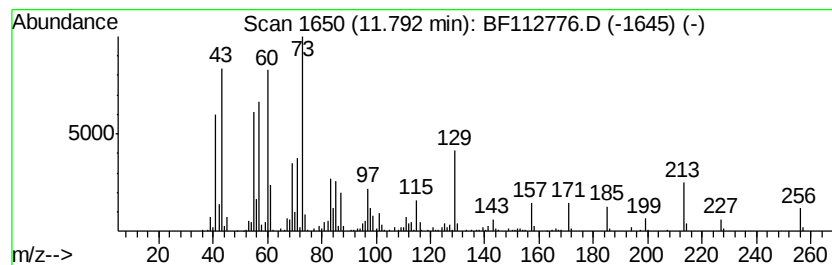
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 BNA_F
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.79	10.86 ng	719131	Phenanthrene-d10		11.25	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5		Dodecanoic acid	200	C12H24O2	000143-07-7	68



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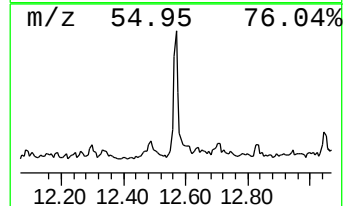
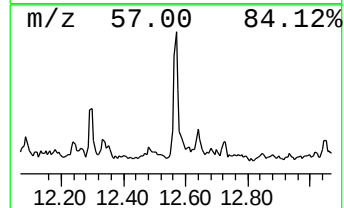
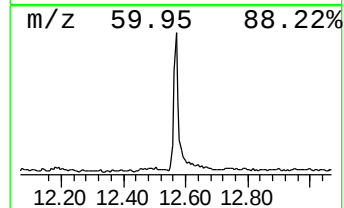
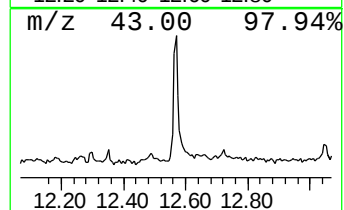
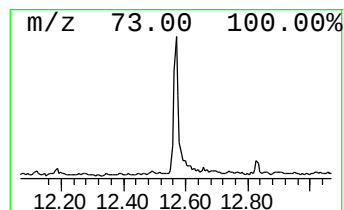
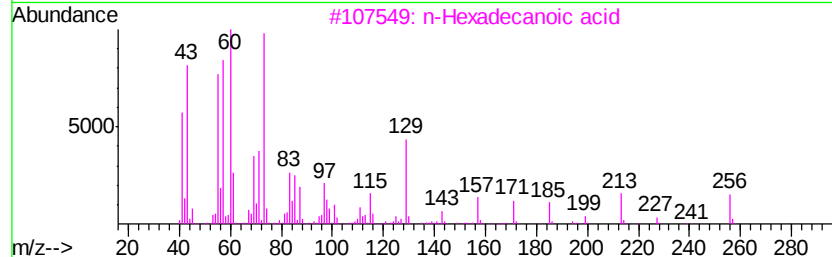
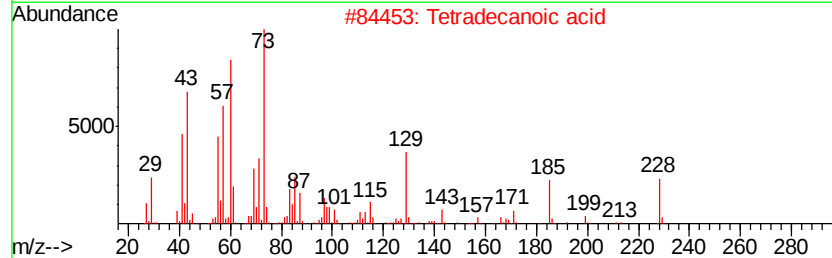
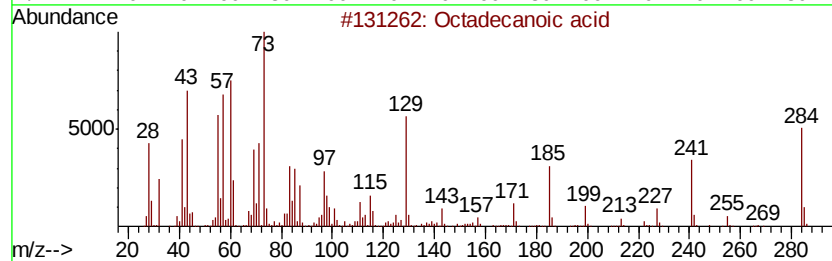
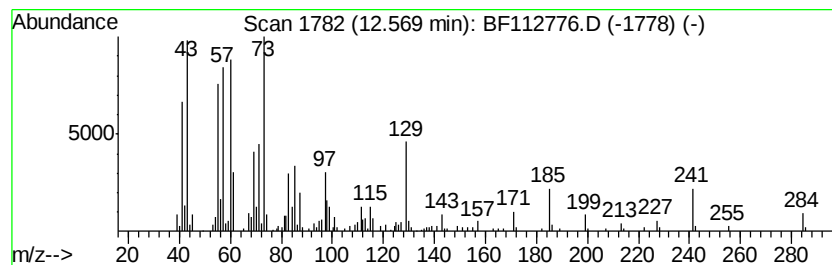
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Peak Number 5 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	3.98 ng	284580	Chrysene-d12	13.87

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
4		Tridecanoic acid	214	C13H26O2	000638-53-9	76
5		n-Decanoic acid	172	C10H20O2	000334-48-5	70



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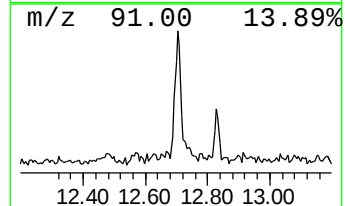
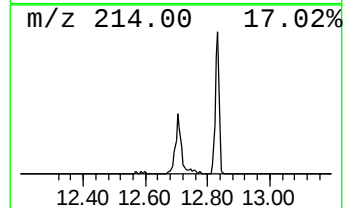
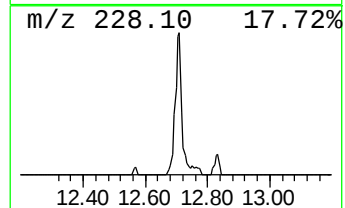
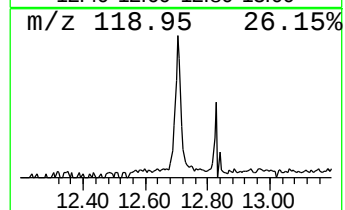
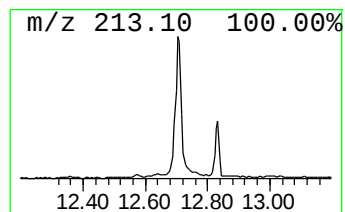
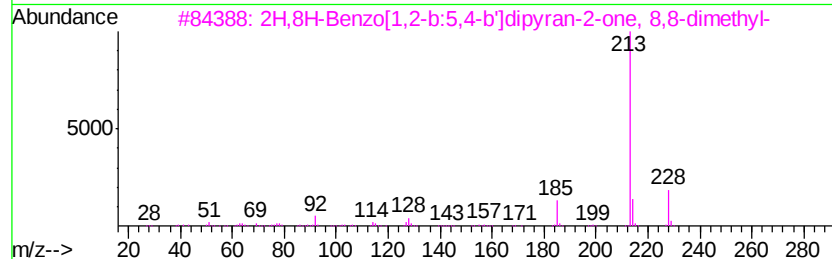
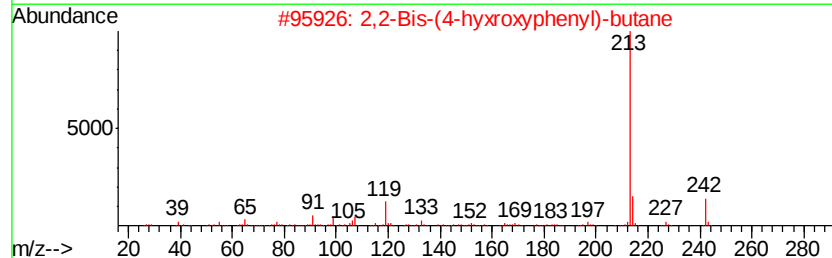
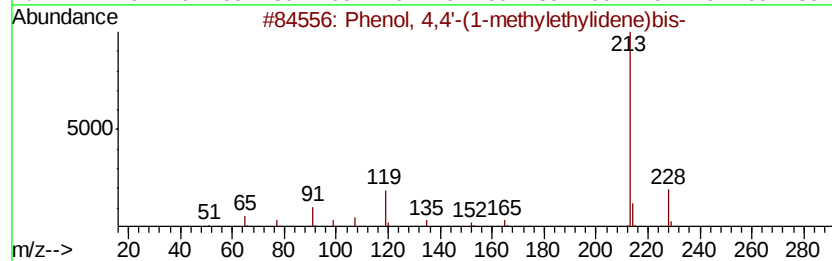
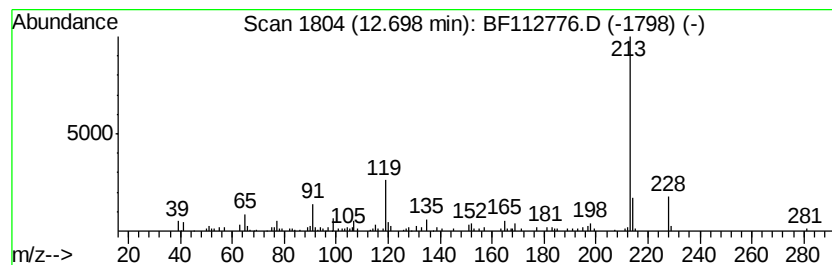
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Peak Number 6 Phenol, 4,4'-(1-methylethyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	2.71 ng	193562	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-(1-methylethylidene)...	228	C15H16O2	000080-05-7	97
2		2,2-Bis-(4-hydroxyphenyl)-butane	242	C16H18O2	000077-40-7	72
3		2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	53
4		2,5-bis(Trimethylsilyl)thiophene	228	C10H20SSi2	017906-71-7	53
5		Indole,6-methyl-2-(2-thienyl)-	213	C13H11NS	296245-70-0	50



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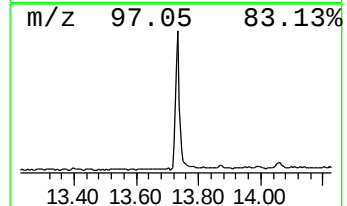
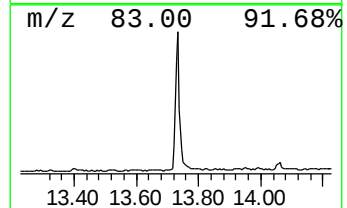
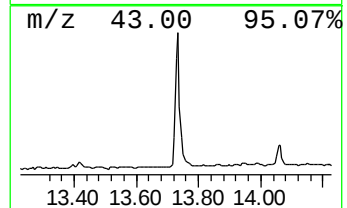
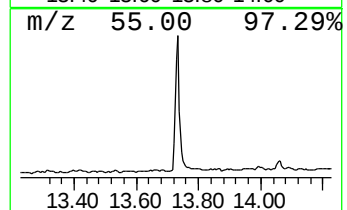
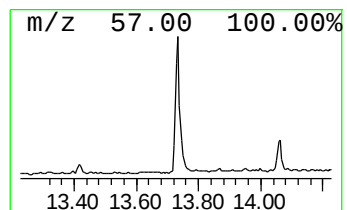
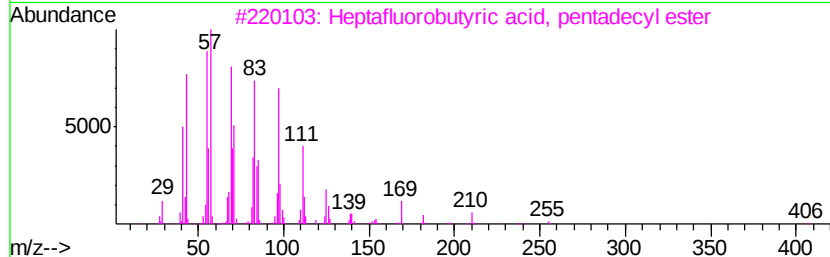
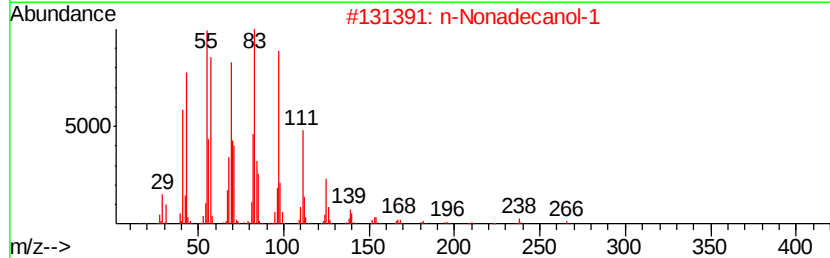
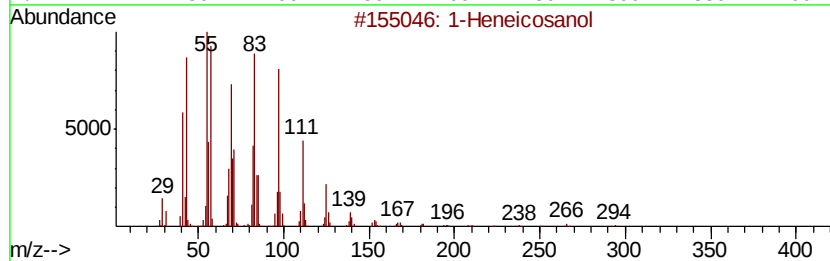
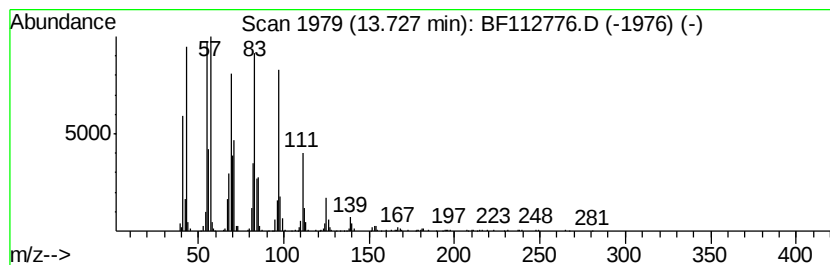
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	11.01 ng	787691	Chrysene-d12	13.87

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol		312	C21H44O	015594-90-8	95
2	n-Nonadecanol-1		284	C19H40O	001454-84-8	95
3	Heptafluorobutyric acid, pentade...		424	C19H31F7O2	959261-23-5	94
4	Pentafluoropropionic acid, penta...		374	C18H31F5O2	959092-08-1	94
5	Behenic alcohol		326	C22H46O	000661-19-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
Data File : BF112776.D
Acq On : 28 Feb 2019 22:12
Operator : JU/SJ
Sample : K1645-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SR-9-D

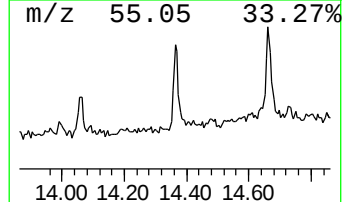
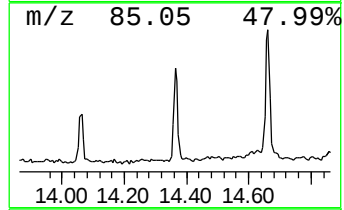
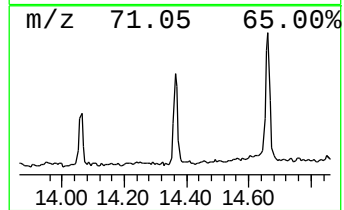
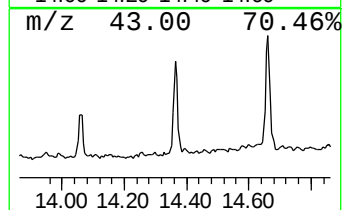
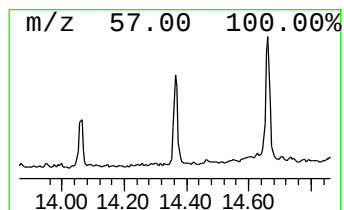
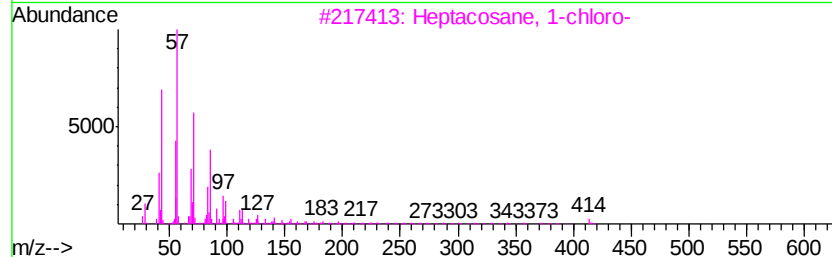
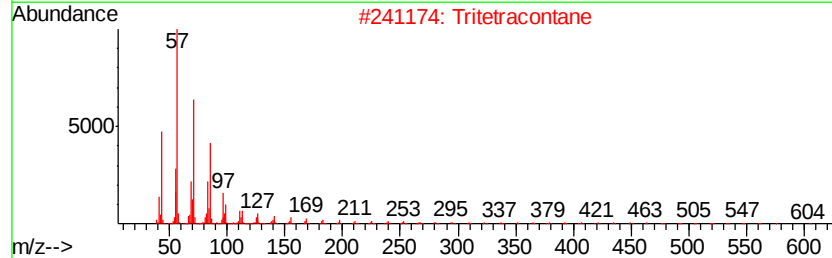
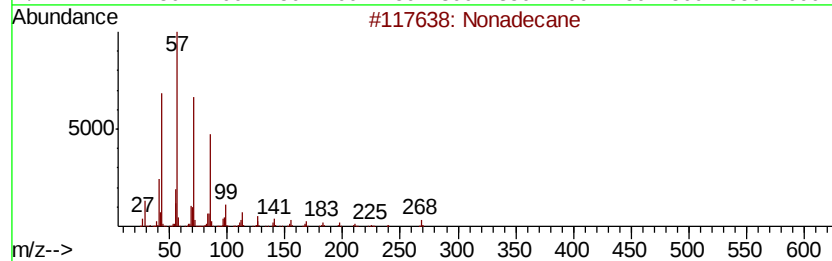
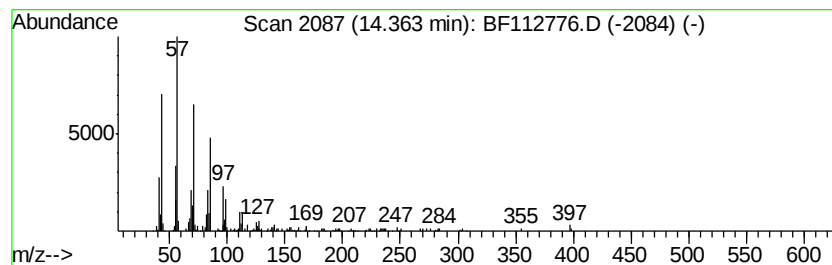
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.36	2.47 ng	177012	Chrysene-d12	13.87

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	93
2	Tritetracontane		605	C43H88	007098-21-7	87
3	Heptacosane, 1-chloro-		414	C27H55Cl	062016-79-9	87
4	Sulfurous acid, butyl dodecyl ester		306	C16H34O3S	1000309-17-9	87
5	Tetratetracontane		619	C44H90	007098-22-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
Data File : BF112776.D
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Sample : K1645-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

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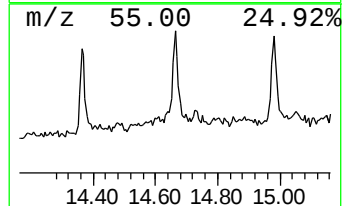
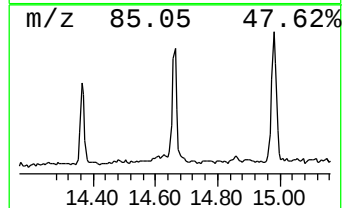
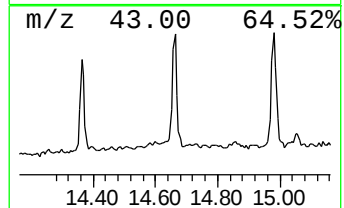
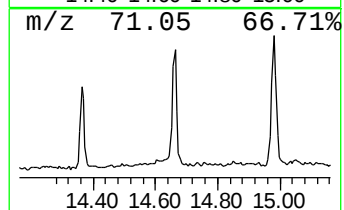
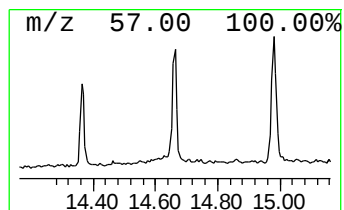
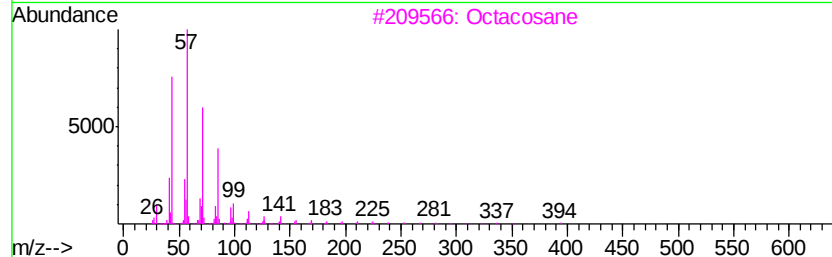
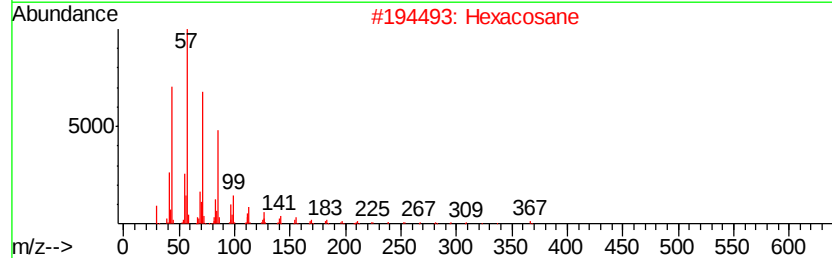
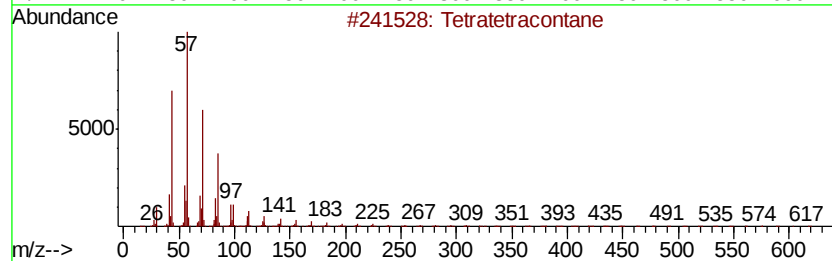
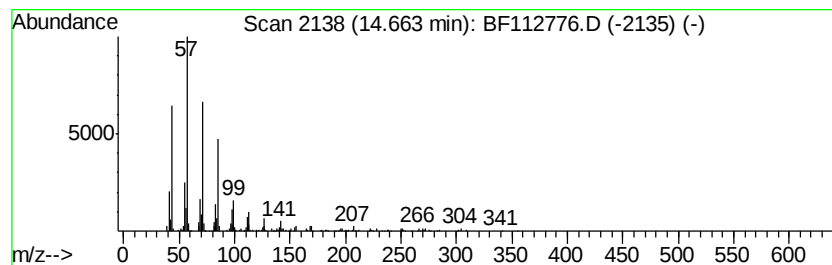
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetratetracontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	2.82 ng	204751	Perylene-d12	15.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetratetracontane	619	C44H90	007098-22-8	90
2		Hexacosane	366	C26H54	000630-01-3	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Heptadecane	240	C17H36	000629-78-7	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
Data File : BF112776.D
Acq On : 28 Feb 2019 22:12
Operator : JU/SJ
Sample : K1645-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

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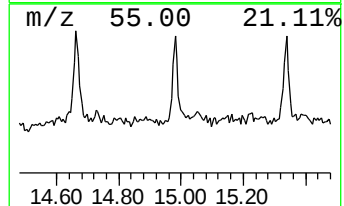
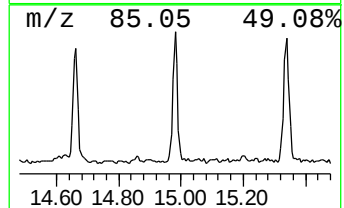
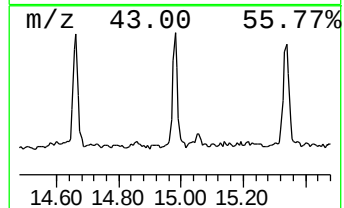
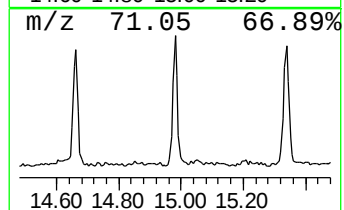
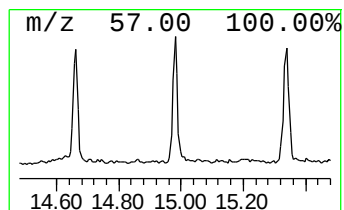
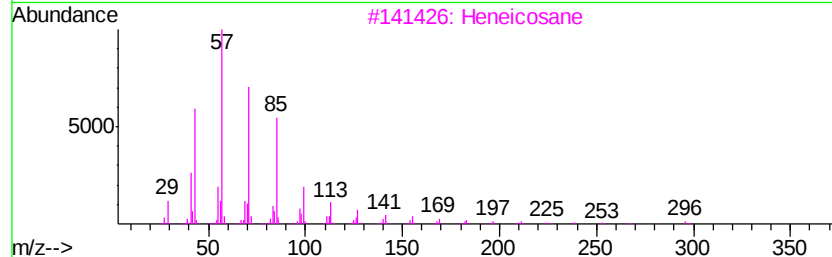
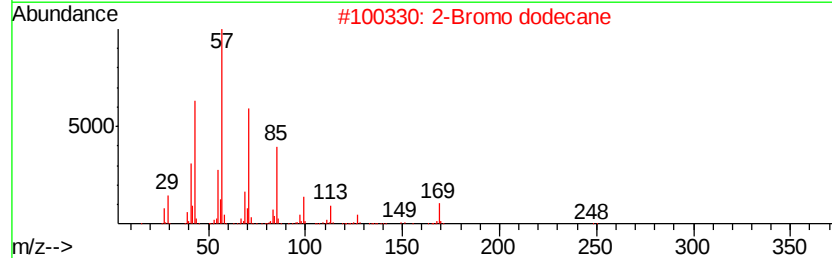
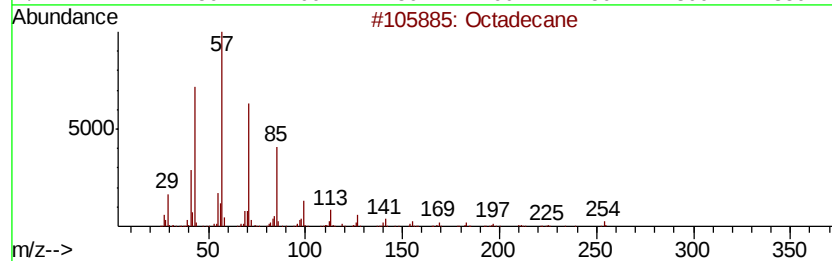
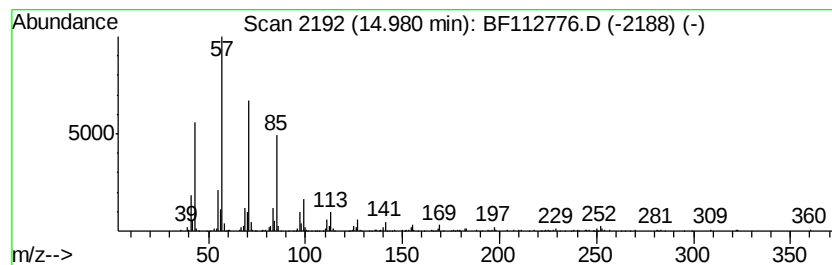
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	3.47 ng	251812	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	95
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
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Acq On : 28 Feb 2019 22:12
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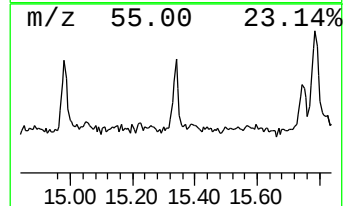
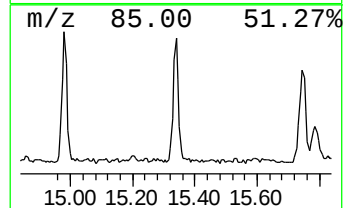
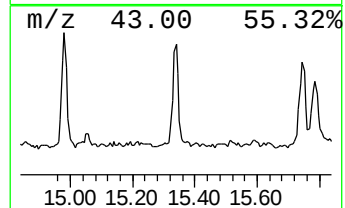
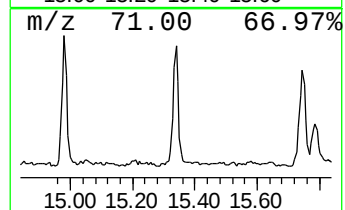
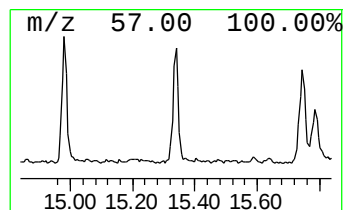
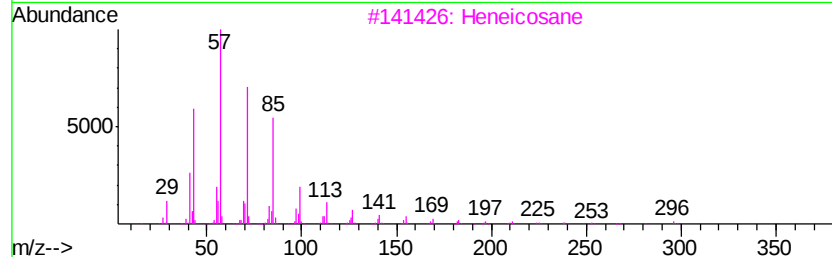
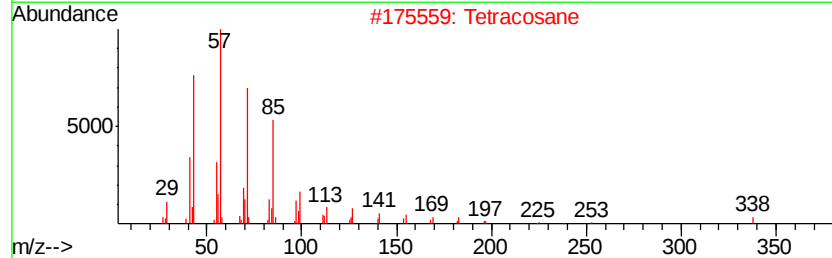
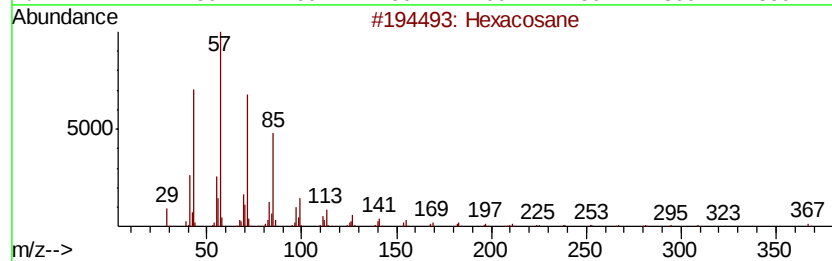
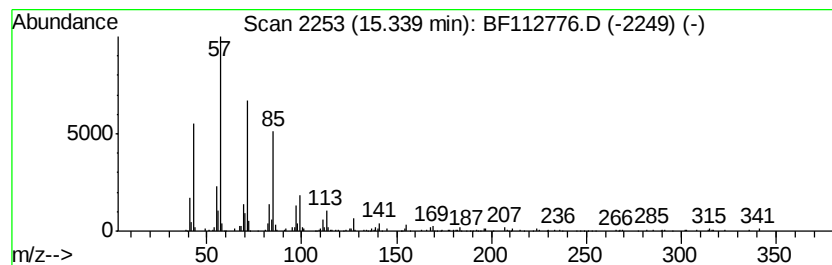
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Hexacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	3.24 ng	235436	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
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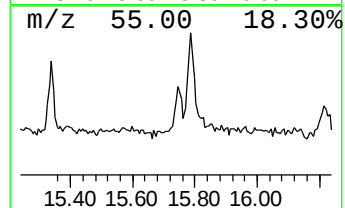
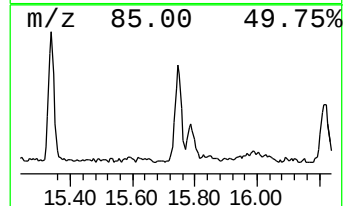
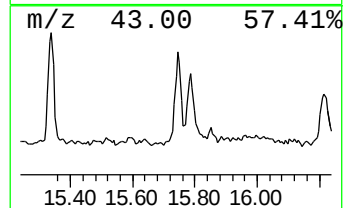
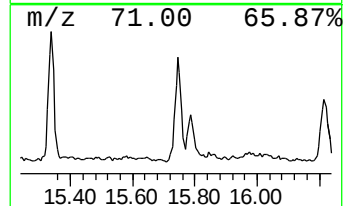
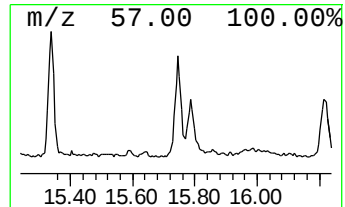
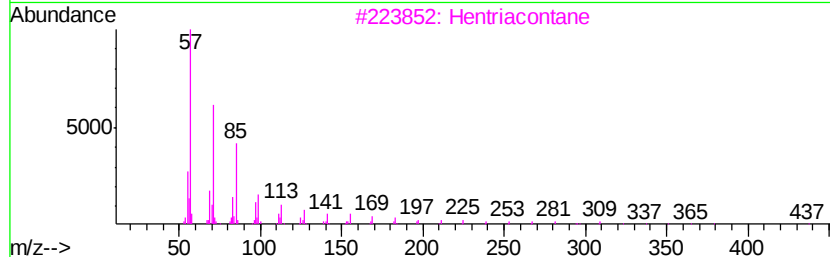
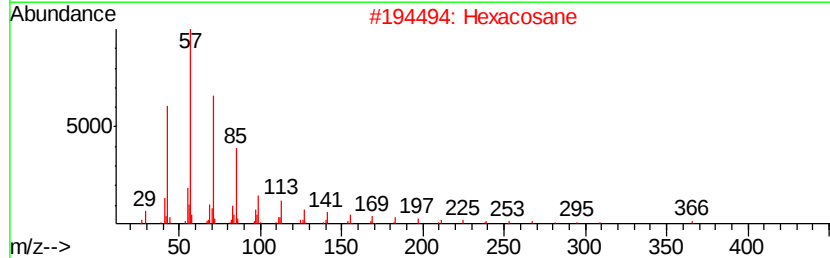
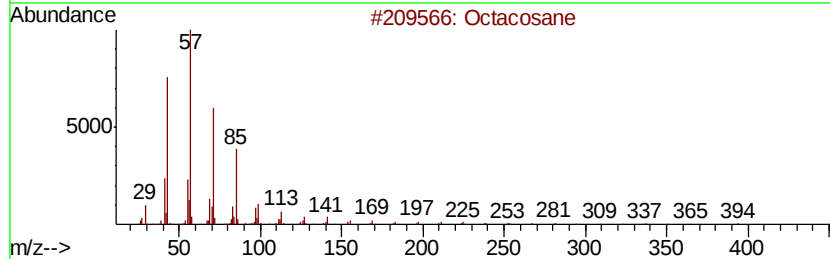
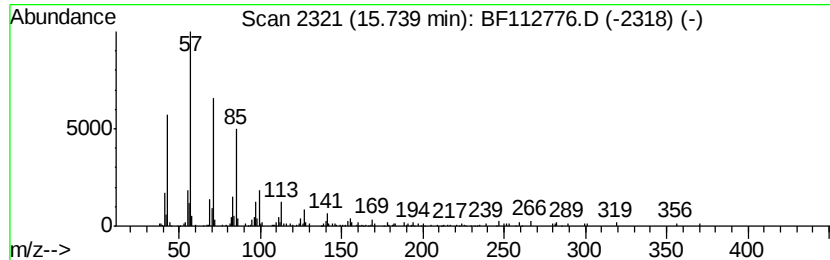
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	2.88 ng	208989	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	87
2		Hexacosane	366	C26H54	000630-01-3	87
3		Hentriacontane	437	C31H64	000630-04-6	86
4		2-Bromotetradecane	276	C14H29Br	074036-95-6	86
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
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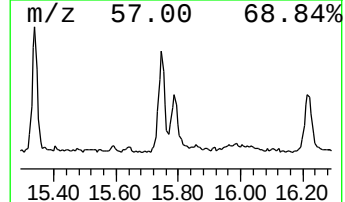
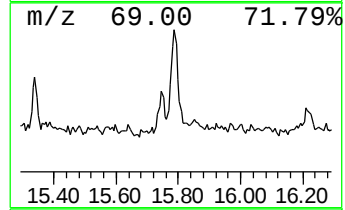
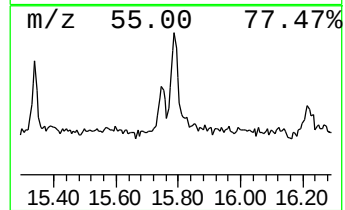
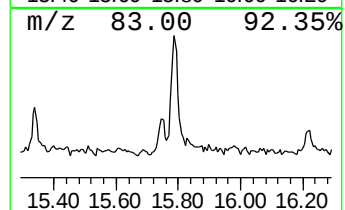
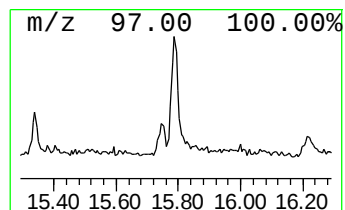
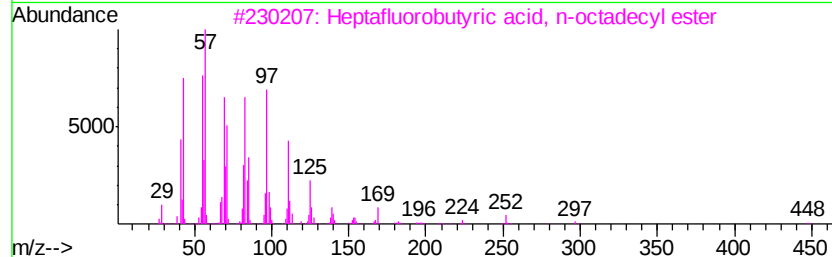
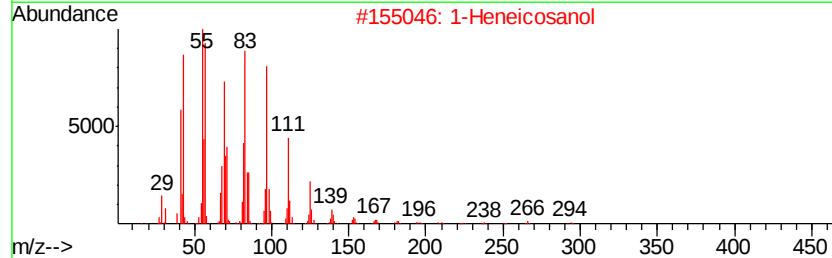
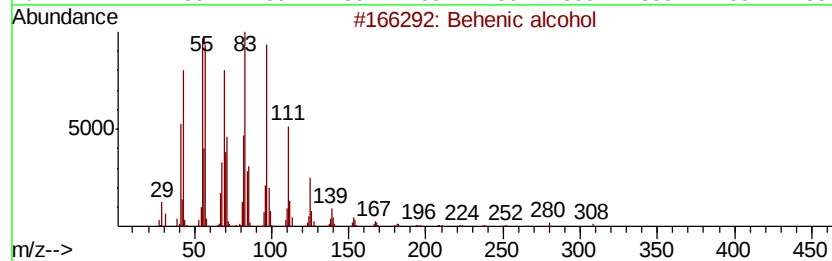
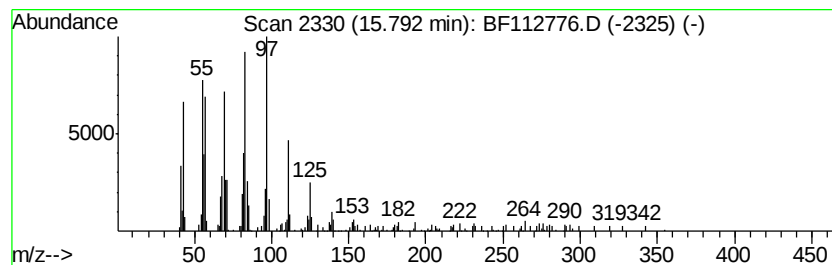
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Behenic alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.79	4.10 ng	297612	Perylene-d12	15.29	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	91
3	Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	91
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5	1-Heptacosanol	396	C27H56O	002004-39-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022819\
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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
3-Penten-2-one, 4...	4.30	2.2	ng	138793	1	6.75	1238220	20.0
unknown6.52	6.52	70.5	ng	4366850	1	6.75	1238220	20.0
n-Hexadecanoic acid	11.79	10.9	ng	719131	4	11.25	1324670	20.0
Octadecanoic acid	12.57	4.0	ng	284580	5	13.87	1430810	20.0
Phenol, 4,4'-(1-m...	12.70	2.7	ng	193562	5	13.87	1430810	20.0
1-Heneicosanol	13.73	11.0	ng	787691	5	13.87	1430810	20.0
Nonadecane	14.36	2.5	ng	177012	5	13.87	1430810	20.0
Tetratetracontane	14.66	2.8	ng	204751	6	15.29	1453130	20.0
Octadecane	14.98	3.5	ng	251812	6	15.29	1453130	20.0
Hexacosane	15.34	3.2	ng	235436	6	15.29	1453130	20.0
Octacosane	15.74	2.9	ng	208989	6	15.29	1453130	20.0
Behenic alcohol	15.79	4.1	ng	297612	6	15.29	1453130	20.0