

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF030719\
 Data File : BF112932.D
 Acq On : 7 Mar 2019 20:12
 Operator : JU/SJ
 Sample : K1760-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Z2-9

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	5.340	548	553	556	rBV	3308982	3095488	78.76%	7.395%
2	6.363	722	727	740	rBV	3266677	3396144	86.41%	8.113%
3	6.498	745	750	753	rBV	3667574	3376635	85.91%	8.066%
4	6.728	785	789	793	rBV	1384142	1115287	28.38%	2.664%
5	6.887	812	816	819	rBV	3138756	2642959	67.24%	6.314%
6	7.287	879	884	887	rBV	2289024	2056253	52.32%	4.912%
7	8.010	1003	1007	1014	rBV	1454368	1373542	34.95%	3.281%
8	9.086	1186	1190	1204	rBV	4233194	3930424	100.00%	9.389%
9	9.210	1208	1211	1221	rVB	232970	230887	5.87%	0.552%
10	9.475	1253	1256	1263	rBV	215459	223275	5.68%	0.533%
11	9.757	1300	1304	1320	rBV	1652151	1668947	42.46%	3.987%
12	10.351	1402	1405	1410	rBV	40610	46351	1.18%	0.111%
13	10.545	1433	1438	1449	rBV	2659770	2732692	69.53%	6.528%
14	11.122	1528	1536	1539	rBV4	26779	51776	1.32%	0.124%
15	11.233	1552	1555	1558	rBV	1768385	1739414	44.26%	4.155%
16	11.292	1563	1565	1570	rVB3	39460	52192	1.33%	0.125%
17	11.710	1631	1636	1639	rBV2	44421	71914	1.83%	0.172%
18	11.775	1643	1647	1656	rVV	792983	874114	22.24%	2.088%
19	12.280	1730	1733	1737	rVB2	90696	82407	2.10%	0.197%
20	12.439	1757	1760	1765	rBV	320972	315554	8.03%	0.754%
21	12.486	1766	1768	1775	rVB5	43021	56979	1.45%	0.136%
22	12.557	1777	1780	1789	rBV2	268422	358115	9.11%	0.855%
23	12.669	1794	1799	1803	rVV	255578	282914	7.20%	0.676%
24	12.816	1820	1824	1828	rBV	3795091	3658195	93.07%	8.739%
25	13.063	1863	1866	1869	rVB	62304	58373	1.49%	0.139%
26	13.098	1869	1872	1881	rVB4	34699	59455	1.51%	0.142%
27	13.404	1921	1924	1928	rVB	69566	68227	1.74%	0.163%
28	13.504	1938	1941	1945	rVB3	34691	46519	1.18%	0.111%
29	13.727	1975	1979	1989	rVB	490328	682439	17.36%	1.630%
30	13.857	1996	2001	2005	rBV	1465252	1628238	41.43%	3.890%
31	14.045	2029	2033	2036	rBV	273965	271612	6.91%	0.649%
32	14.345	2081	2084	2090	rBV	507507	494567	12.58%	1.181%
33	14.639	2131	2134	2142	rBV	605069	604814	15.39%	1.445%
34	14.710	2143	2146	2149	rVB	220848	203531	5.18%	0.486%

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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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
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35	14.868	2169	2173	2182	rVB	117450	251226	6.39%	0.600%
36	14.957	2184	2188	2193	rVV	596581	646720	16.45%	1.545%
37	15.145	2217	2220	2223	rBV	78517	83385	2.12%	0.199%
38	15.198	2227	2229	2235	rVV	76858	102791	2.62%	0.246%
39	15.263	2235	2240	2244	rVV	1148814	1414911	36.00%	3.380%
40	15.310	2244	2248	2256	rVB	531604	701032	17.84%	1.675%
41	15.715	2312	2317	2322	rBV	387282	574641	14.62%	1.373%
42	16.180	2392	2396	2401	rBV	212493	353086	8.98%	0.843%
43	16.727	2485	2489	2496	rVB2	92597	182638	4.65%	0.436%

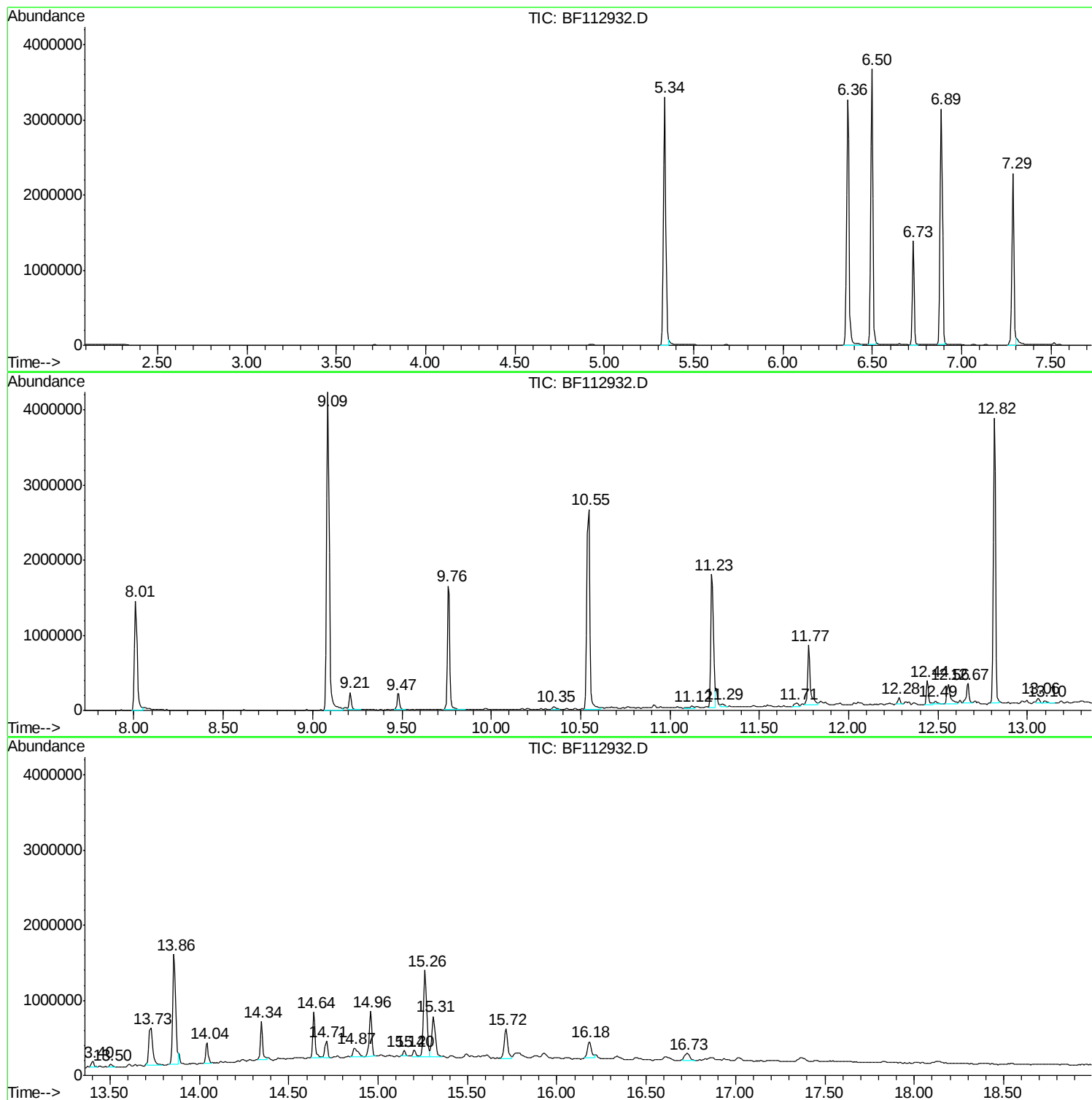
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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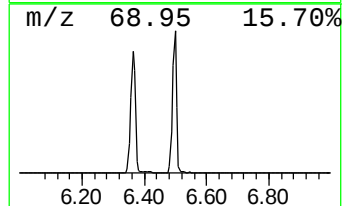
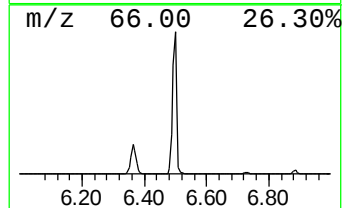
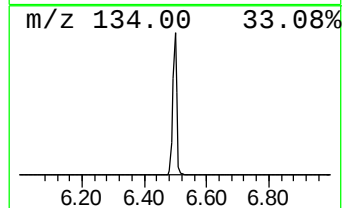
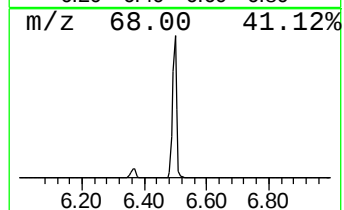
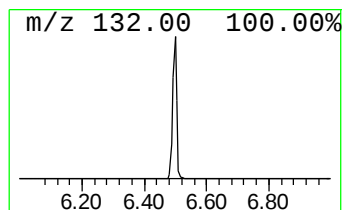
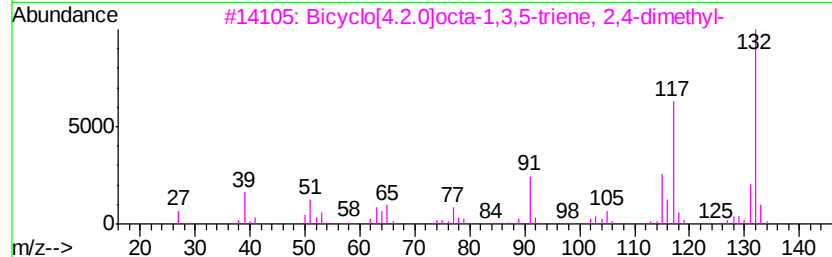
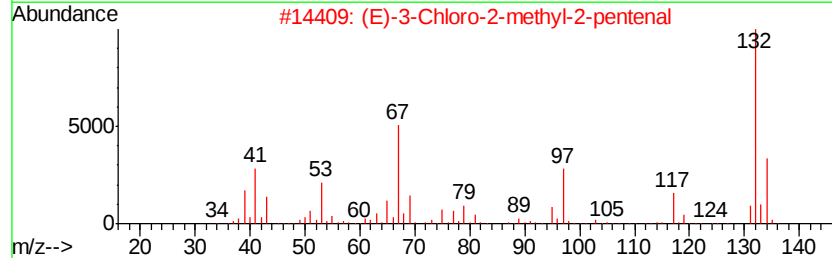
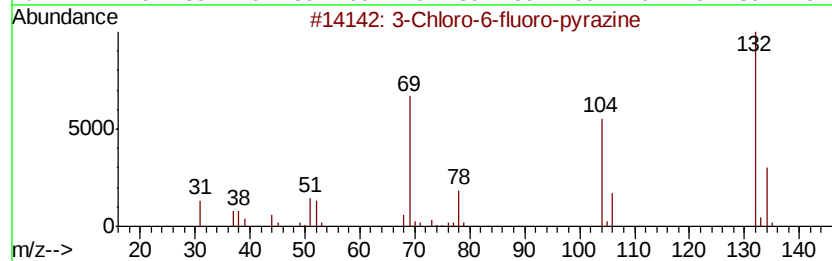
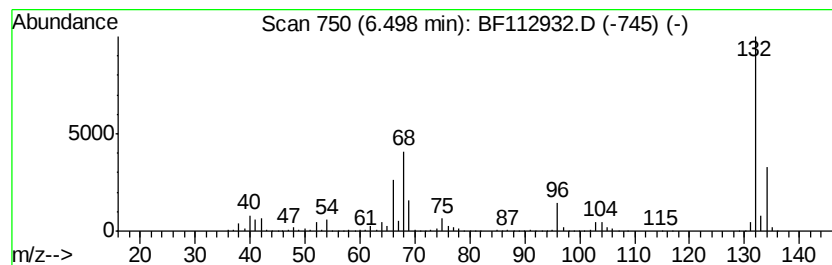
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	60.55 ng	3376640	1,4-Dichlorobenzene-d4	6.73

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	16
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Xenon	132	Xe	007440-63-3	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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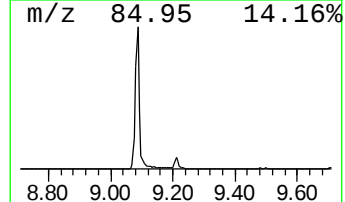
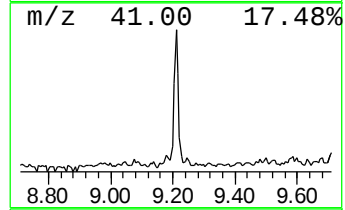
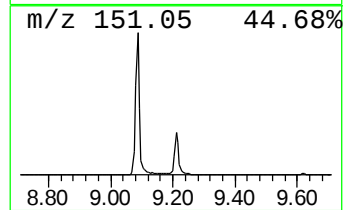
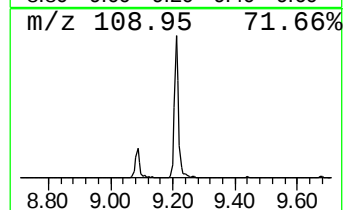
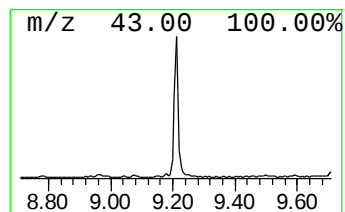
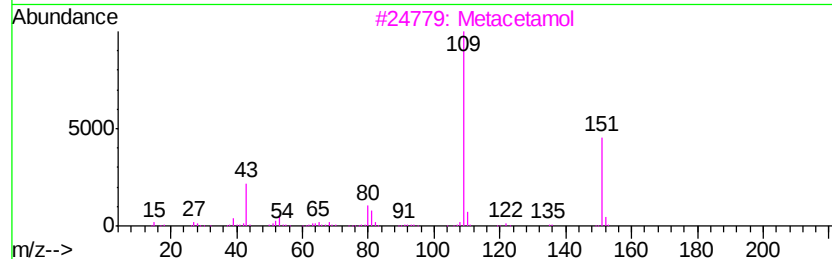
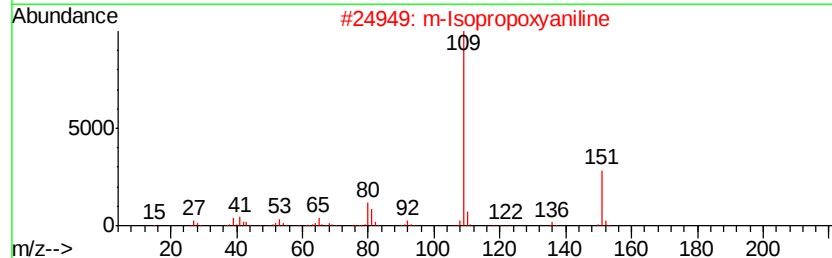
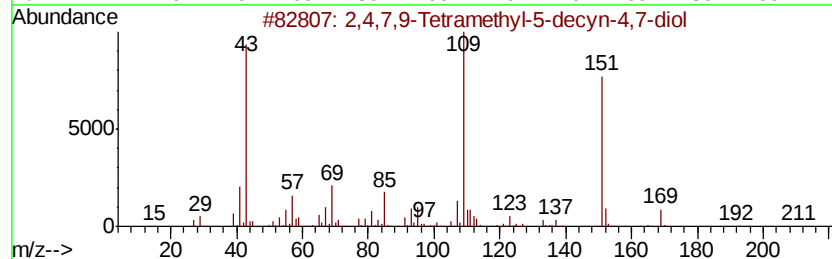
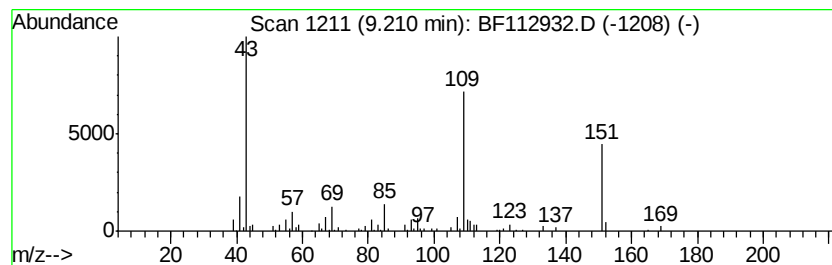
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	2.77 ng	230887	Acenaphthene-d10	9.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90
2		m-Isopropoxyvaniline	151	C9H13NO	041406-00-2	53
3		Metacetamol	151	C8H9NO2	000621-42-1	53
4		Acetaminophen	151	C8H9NO2	000103-90-2	42
5		Butanamide, N-acetyl-N-(4-hydrox...	221	C12H15NO3	1000107-08-7	42



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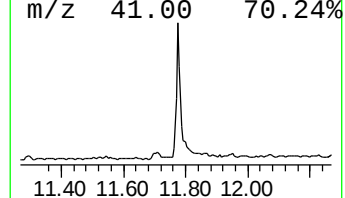
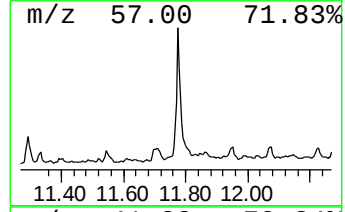
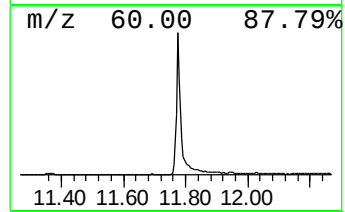
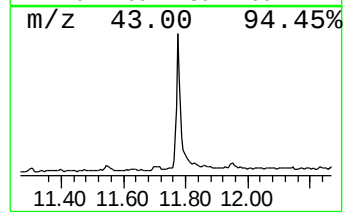
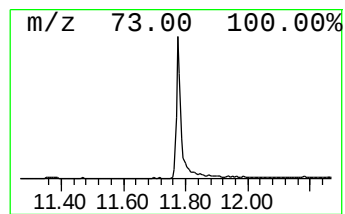
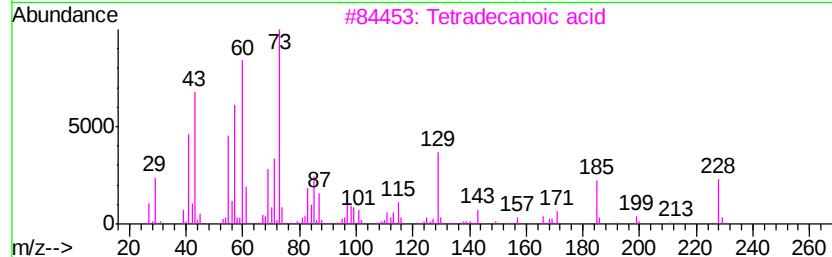
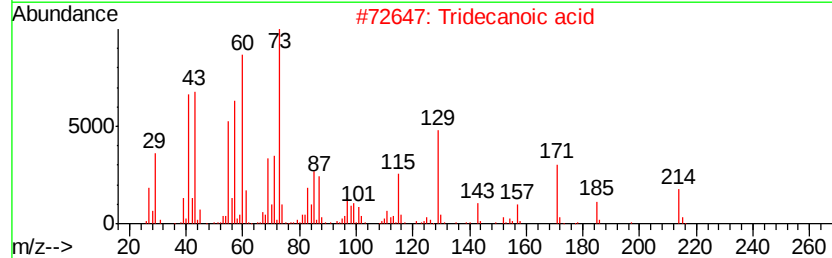
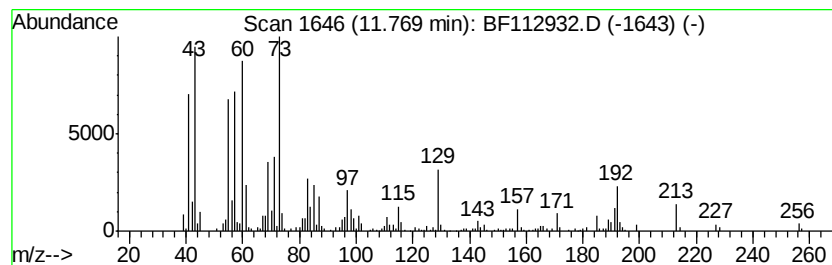
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	10.05 ng	874114	Phenanthrene-d10	11.23

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		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5		Oxalic acid, dodecyl propyl ester	300	C17H32O4	1000309-26-5	30



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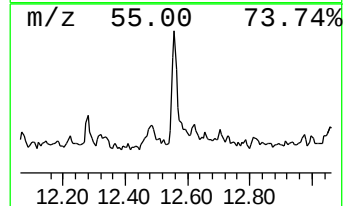
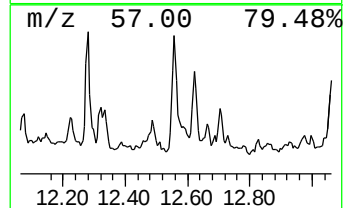
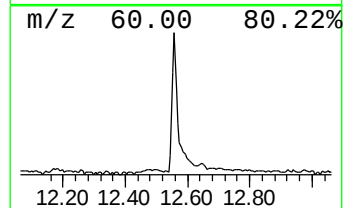
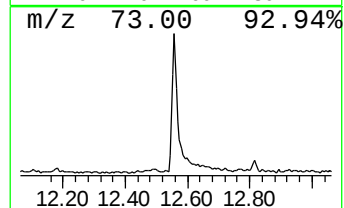
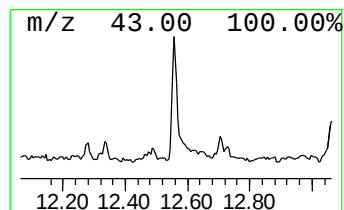
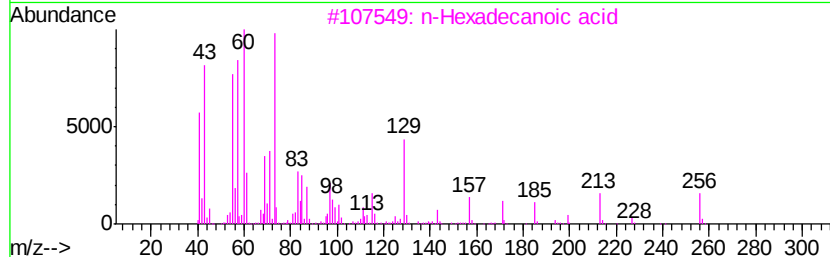
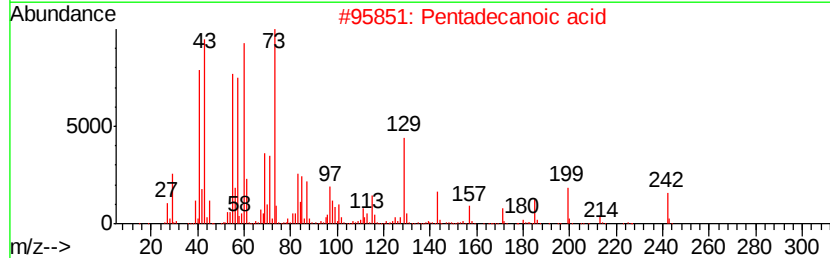
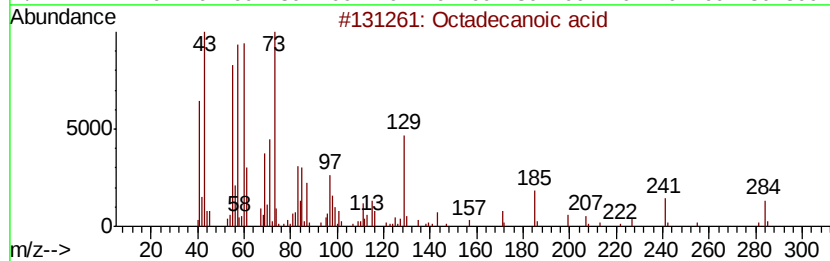
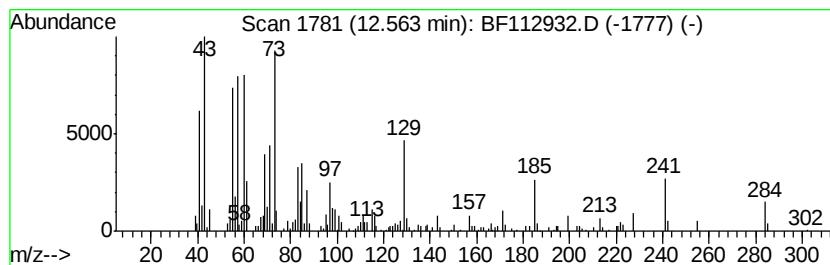
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	4.40 ng	358115	Chrysene-d12	13.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5		Octadecanoic acid, 2-(2-hydroxy...	372	C22H44O4	000106-11-6	72



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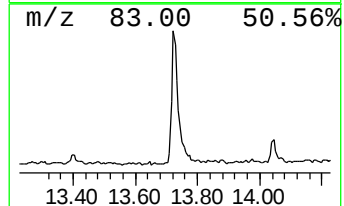
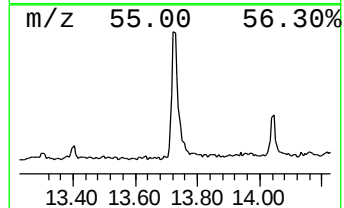
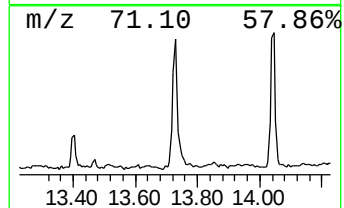
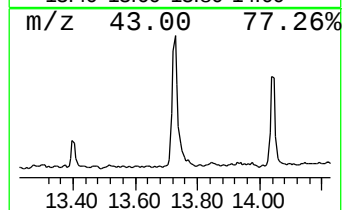
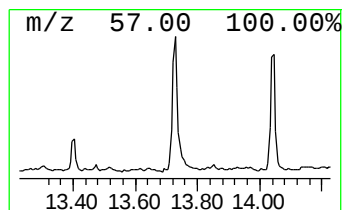
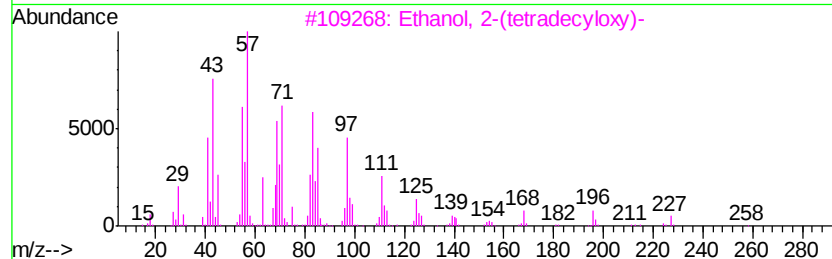
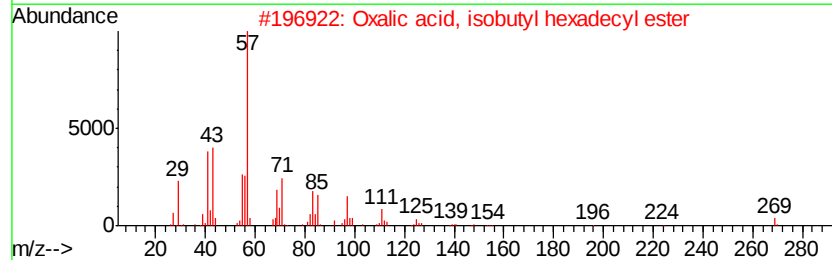
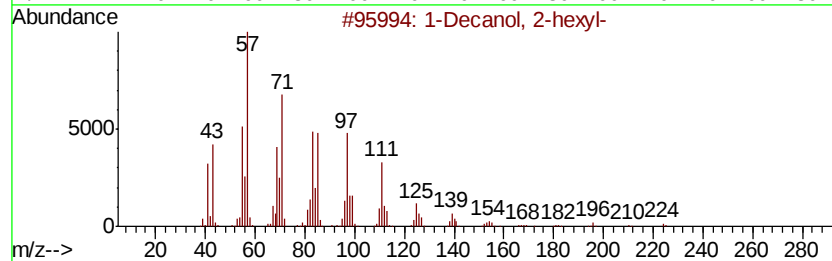
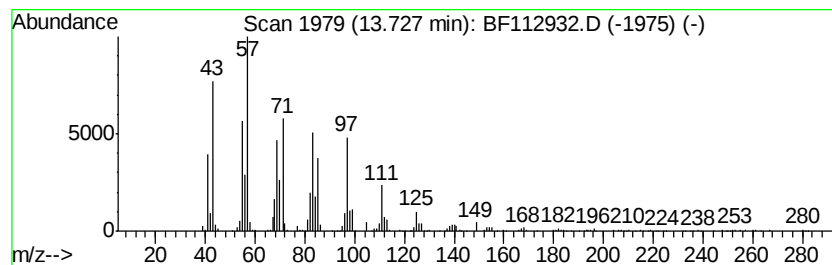
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 6 1-Decanol, 2-hexyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	8.38 ng	682439	Chrysene-d12	13.86
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Decanol, 2-hexyl-	242 C16H34O	002425-77-6 91
2		Oxalic acid, isobutyl hexadecyl ...	370 C22H42O4	1000309-38-1 91
3		Ethanol, 2-(tetradecyl)-	258 C16H34O2	002136-70-1 90
4		Cyclotetradecane	196 C14H28	000295-17-0 87
5		Heptafluorobutyric acid, n-octad...	466 C22H37F7O2	000400-57-7 87



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Operator : JU/SJ
Sample : K1760-07
Misc :
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ClientSampleId :
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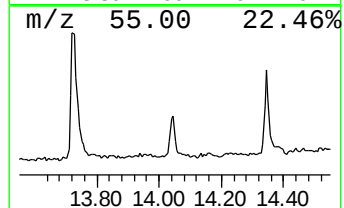
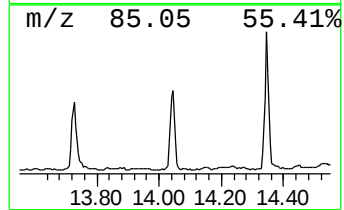
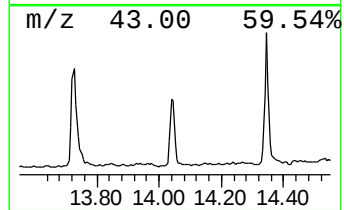
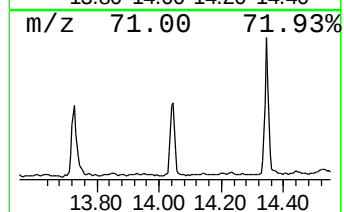
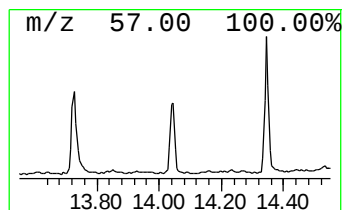
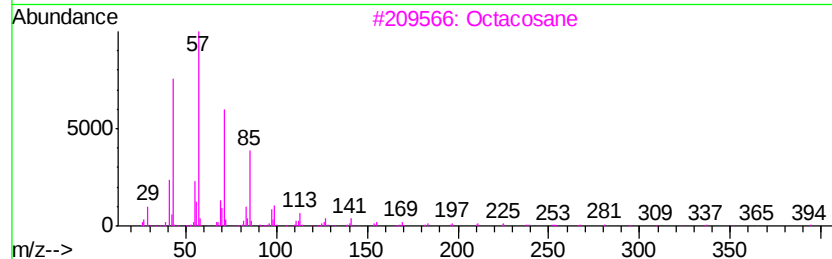
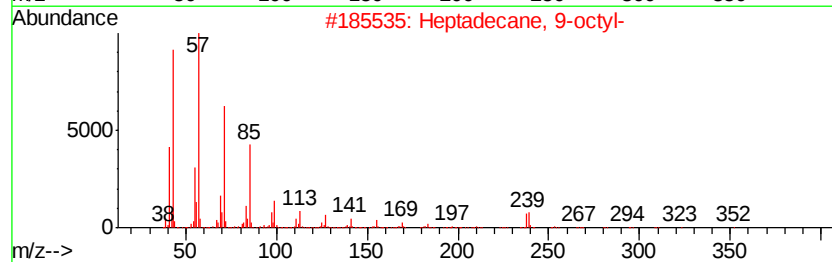
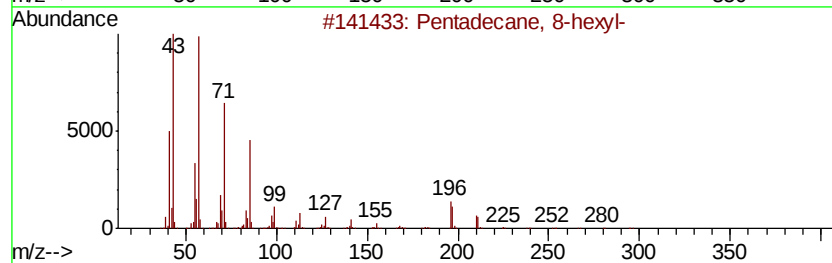
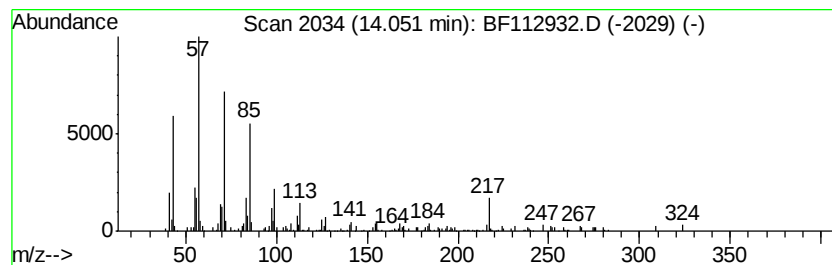
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecane, 8-hexyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.05	3.34 ng	271612	Chrysene-d12	13.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
3		Octacosane	394	C28H58	000630-02-4	87
4		triacontane	422	C30H62	000638-68-6	87
5		Heptadecane	240	C17H36	000629-78-7	87



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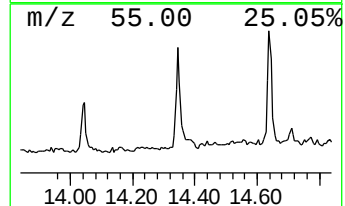
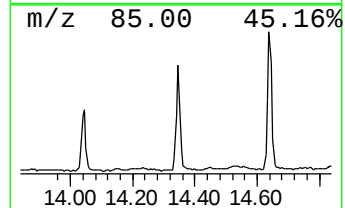
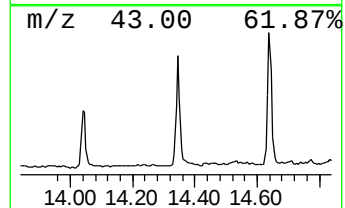
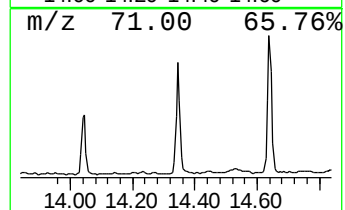
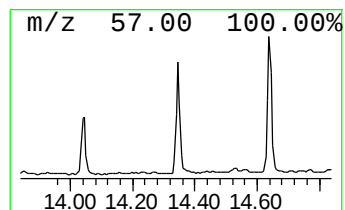
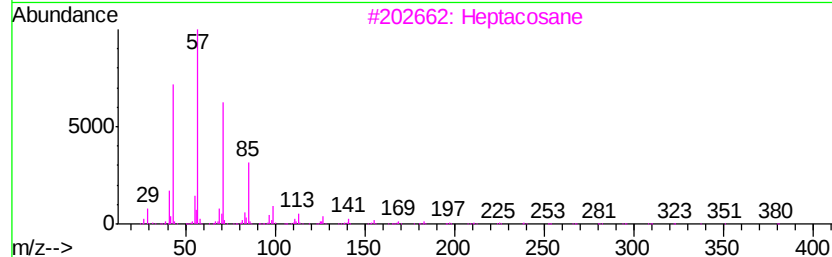
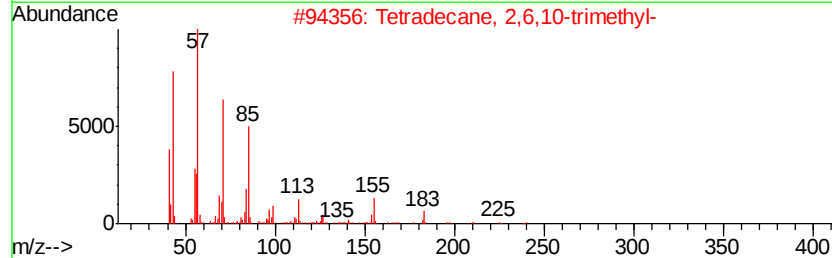
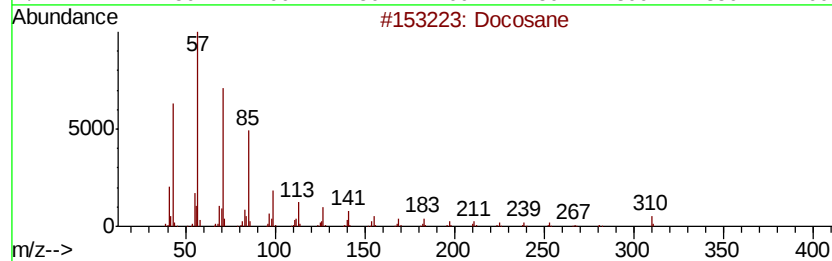
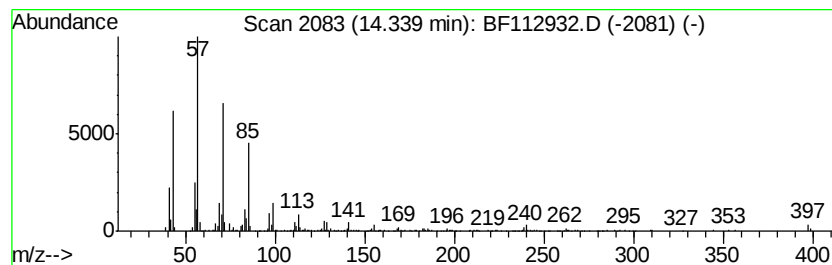
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Docosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.34	6.07 ng	494567	Chrysene-d12	13.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	96
2		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	93
3		Heptacosane	380	C27H56	000593-49-7	91
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030719\
Data File : BF112932.D
Acq On : 7 Mar 2019 20:12
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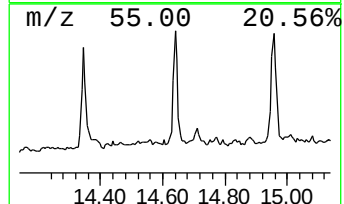
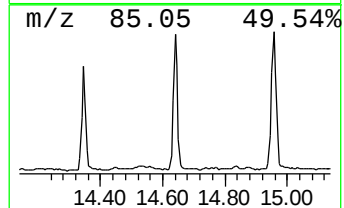
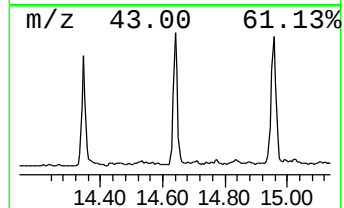
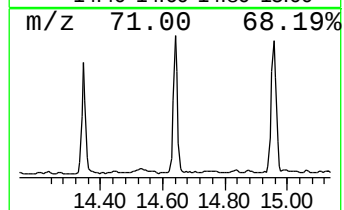
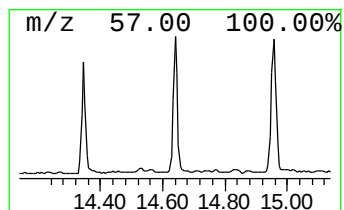
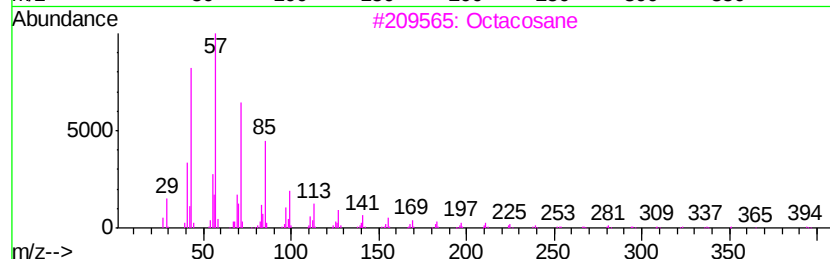
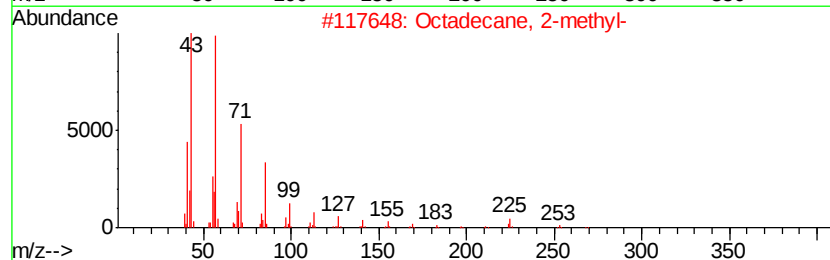
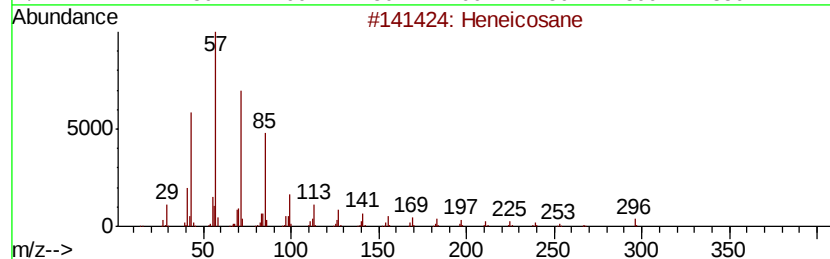
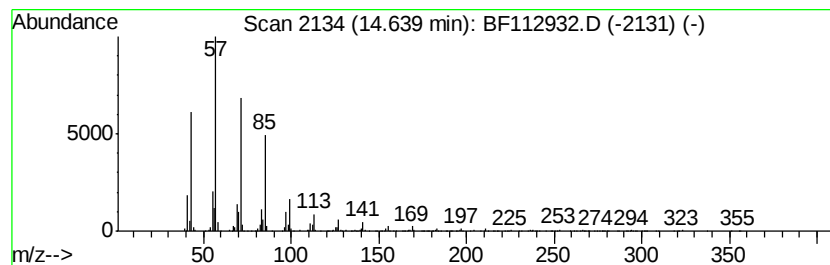
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	8.55 ng	604814	Perylene-d12	15.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	95
2	Octadecane, 2-methyl-		268	C19H40	001560-88-9	93
3	Octacosane		394	C28H58	000630-02-4	91
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
5	Tetracosane		338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030719\
Data File : BF112932.D
Acq On : 7 Mar 2019 20:12
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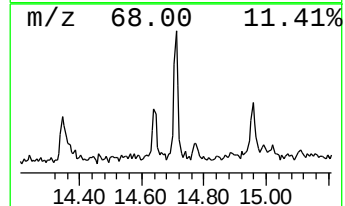
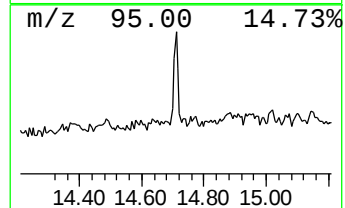
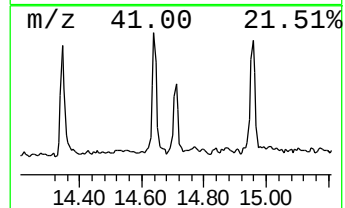
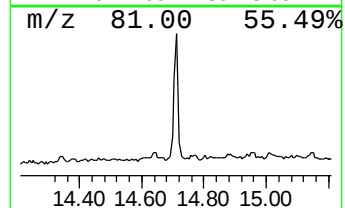
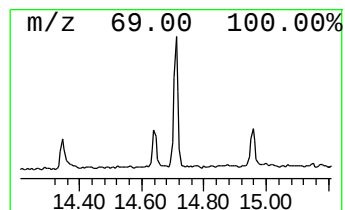
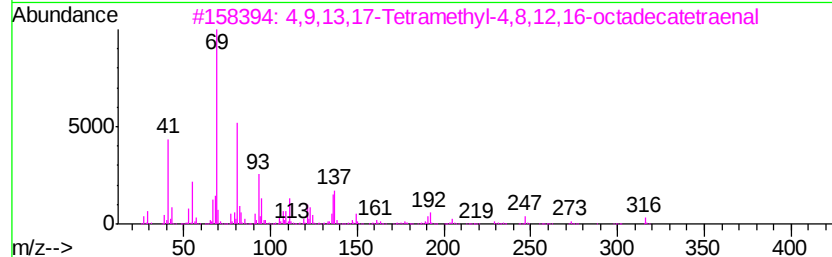
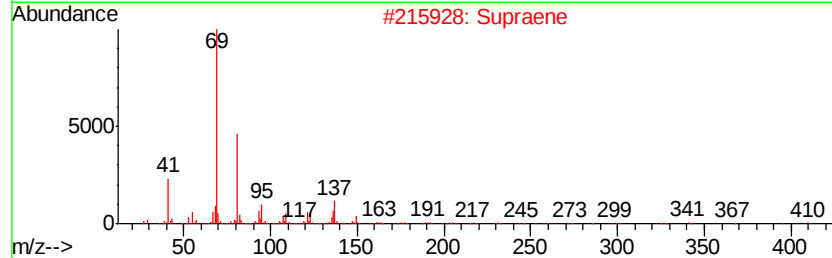
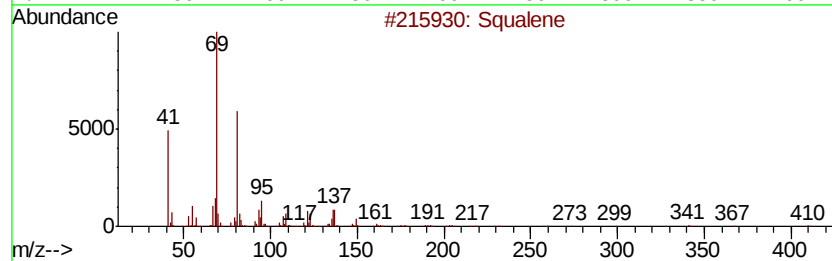
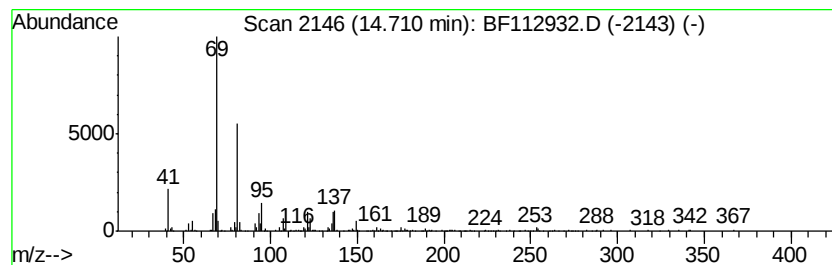
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Squalene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	2.88 ng	203531	Perylene-d12	15.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	91
2		Supraene	410	C30H50	007683-64-9	91
3		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	83
4		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64
5		2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030719\
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Misc :
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Instrument :
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ClientSampleId :
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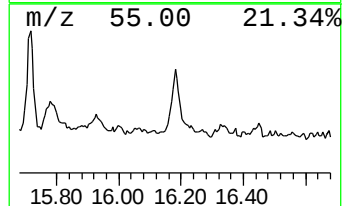
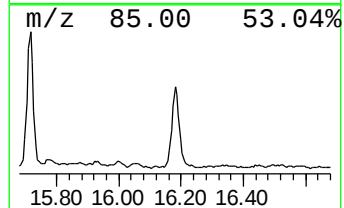
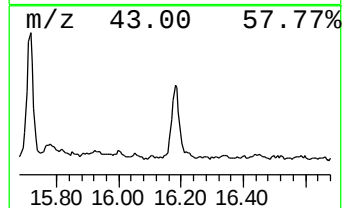
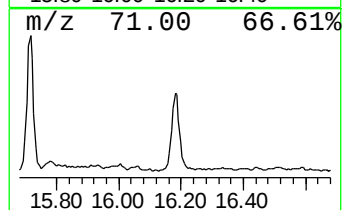
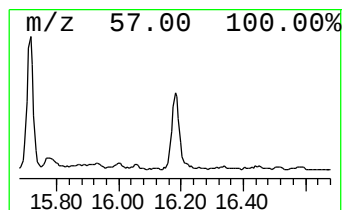
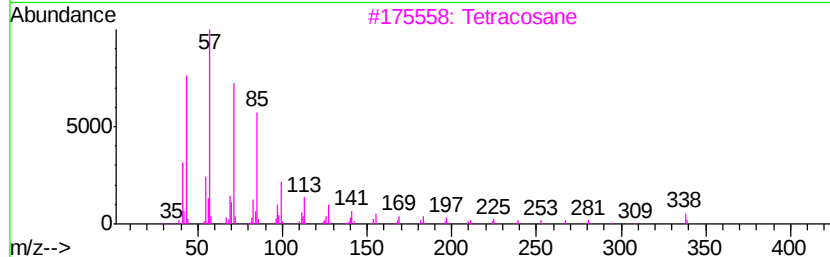
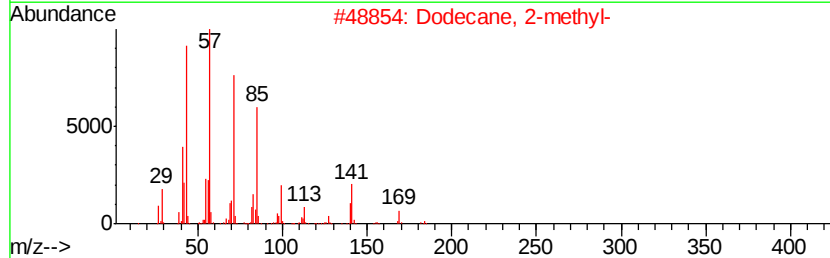
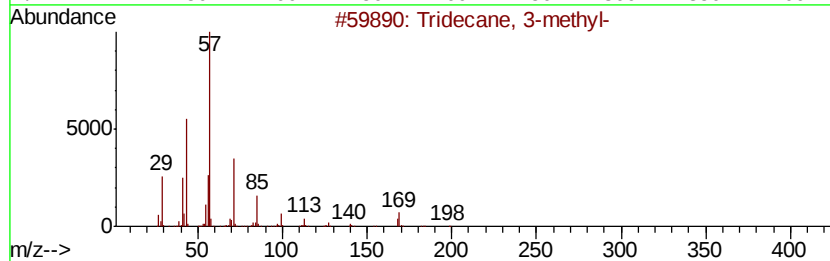
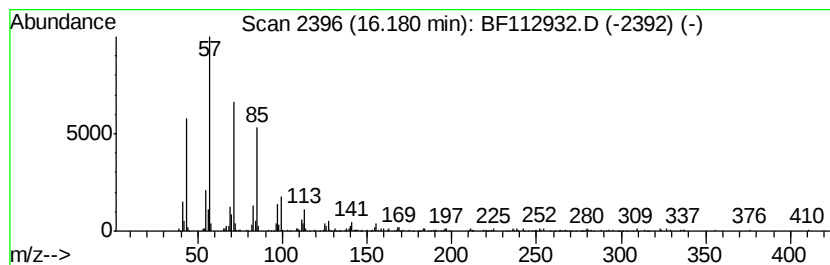
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Tridecane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	4.99 ng	353086	Perylene-d12	15.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 3-methyl-	198	C14H30	006418-41-3	87
2		Dodecane, 2-methyl-	184	C13H28	001560-97-0	86
3		Tetracosane	338	C24H50	000646-31-1	86
4		Heptadecane	240	C17H36	000629-78-7	86
5		Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030719\
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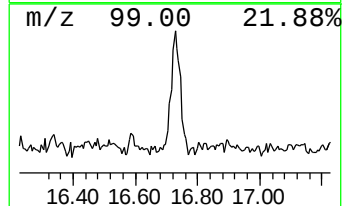
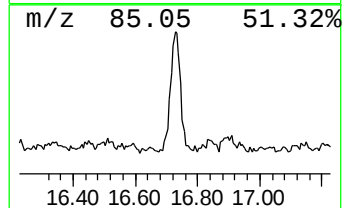
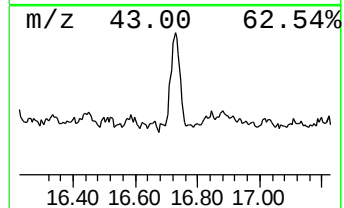
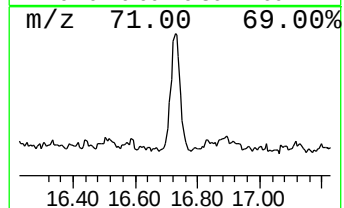
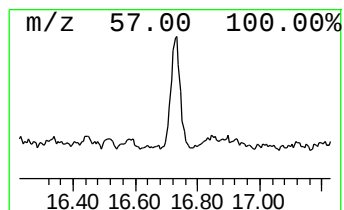
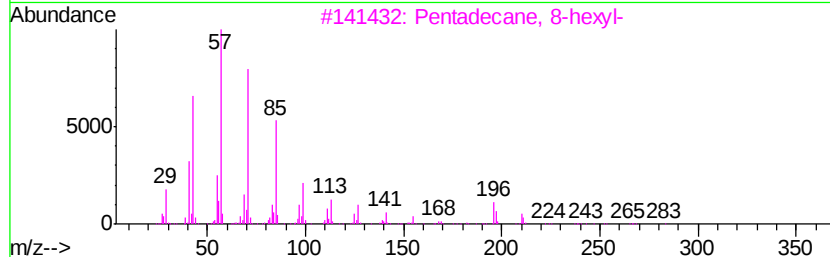
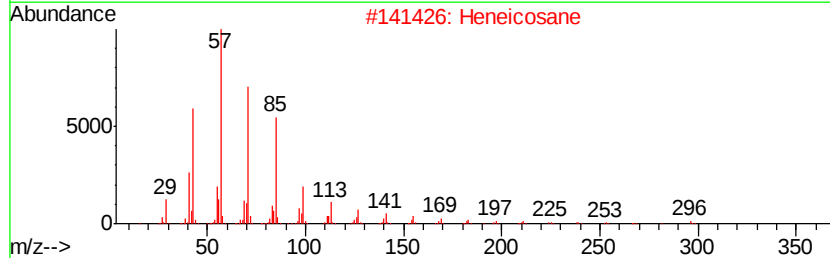
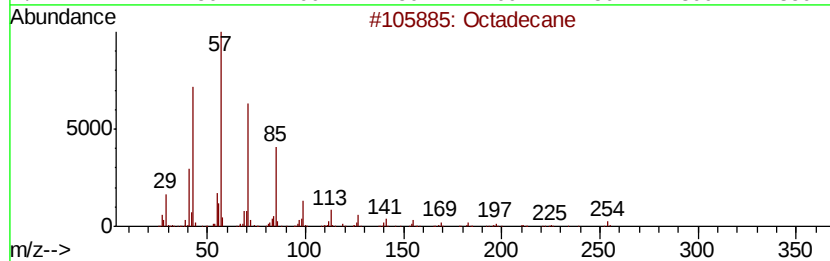
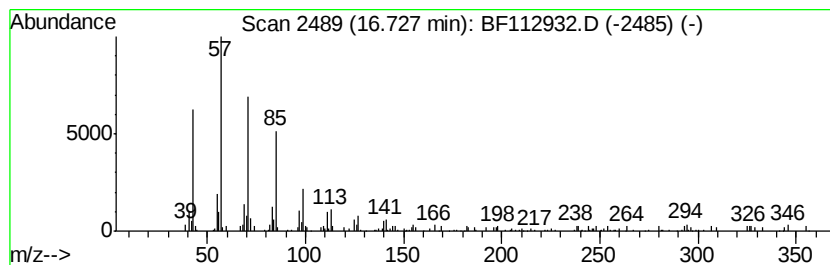
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Octadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.73	2.58 ng	182638	Perylene-d12	15.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	94
2		Heneicosane	296	C21H44	000629-94-7	93
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	86
4		Hexacosane	366	C26H54	000630-01-3	80
5		Docosane	310	C22H46	000629-97-0	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF030719\
 Data File : BF112932.D
 Acq On : 7 Mar 2019 20:12
 Operator : JU/SJ
 Sample : K1760-07
 Misc :
 ALS Vial : 17 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

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					#	RT	Resp	Conc
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2,4,7,9-Tetrameth...	9.21	2.8	ng	230887	3	9.76	1668950	20.0
n-Hexadecanoic acid	11.77	10.1	ng	874114	4	11.23	1739410	20.0
Octadecanoic acid	12.56	4.4	ng	358115	5	13.86	1628240	20.0
1-Decanol, 2-hexyl-	13.73	8.4	ng	682439	5	13.86	1628240	20.0
Pentadecane, 8-he...	14.05	3.3	ng	271612	5	13.86	1628240	20.0
Docosane	14.34	6.1	ng	494567	5	13.86	1628240	20.0
Heneicosane	14.64	8.6	ng	604814	6	15.26	1414910	20.0
Squalene	14.71	2.9	ng	203531	6	15.26	1414910	20.0
Tridecane, 3-methyl-	16.18	5.0	ng	353086	6	15.26	1414910	20.0
Octadecane	16.73	2.6	ng	182638	6	15.26	1414910	20.0