

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
 Data File : BF107629.D
 Acq On : 24 Jul 2018 19:52
 Operator : JU/SJ
 Sample : J4098-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GL-TOPSOIL-02

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-BF072018.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	5.207	484	488	508	rBV	531675	771430	15.97%	1.412%
2	5.607	552	556	592	rBV	3370387	4809565	99.54%	8.806%
3	6.607	721	726	745	rBV	3240585	4667365	96.59%	8.546%
4	6.742	745	749	768	rVB	4207426	4831972	100.00%	8.847%
5	6.966	784	787	807	rVB	1249165	1471756	30.46%	2.695%
6	7.125	810	814	829	rBV	3456408	3604884	74.60%	6.600%
7	7.531	879	883	906	rBV	2469973	3025626	62.62%	5.540%
8	8.248	1001	1005	1026	rBV	1652451	1851709	38.32%	3.390%
9	9.325	1182	1188	1196	rBV	4219144	4212053	87.17%	7.712%
10	9.666	1243	1246	1249	rVB2	38698	51699	1.07%	0.095%
11	9.713	1251	1254	1260	rBV	700081	608151	12.59%	1.113%
12	9.872	1278	1281	1285	rBV	46308	53613	1.11%	0.098%
13	10.007	1301	1304	1314	rBV2	1683437	1685539	34.88%	3.086%
14	10.119	1320	1323	1325	rBV2	61097	50910	1.05%	0.093%
15	10.801	1435	1439	1448	rBV	2260435	2462280	50.96%	4.508%
16	10.907	1452	1457	1460	rVB4	62731	92164	1.91%	0.169%
17	11.342	1529	1531	1534	rBV	66622	67170	1.39%	0.123%
18	11.501	1555	1558	1561	rBV	1452430	1500270	31.05%	2.747%
19	11.530	1561	1563	1567	rVV2	333288	333531	6.90%	0.611%
20	11.583	1569	1572	1575	rBV	41997	51177	1.06%	0.094%
21	11.972	1636	1638	1641	rVB	90302	78966	1.63%	0.145%
22	12.013	1641	1645	1648	rBV	516655	541015	11.20%	0.991%
23	12.189	1673	1675	1681	rVB3	193522	199050	4.12%	0.364%
24	12.624	1747	1749	1751	rBV	99936	83667	1.73%	0.153%
25	12.719	1762	1765	1769	rBV	539426	576641	11.93%	1.056%
26	12.748	1769	1770	1773	rVB	142761	118990	2.46%	0.218%
27	12.954	1799	1805	1812	rBV	496830	585788	12.12%	1.073%
28	13.089	1824	1828	1835	rBV	3195097	3056142	63.25%	5.596%
29	13.519	1898	1901	1904	rVB3	277356	253665	5.25%	0.464%
30	13.966	1974	1977	1980	rBV2	561831	596172	12.34%	1.092%
31	14.154	2005	2009	2012	rBV2	1349032	1463756	30.29%	2.680%
32	14.595	2081	2084	2090	rVB2	862809	1030779	21.33%	1.887%
33	14.995	2149	2152	2157	rVB	322616	337744	6.99%	0.618%
34	15.271	2195	2199	2203	rVB	2383244	2837103	58.72%	5.195%

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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	15.689	2267	2270	2276	rVB	599895	811972	16.80%	1.487%
36	16.142	2342	2347	2353	rVB	2023738	3092671	64.00%	5.663%
37	16.436	2394	2397	2409	rVB2	517269	1044885	21.62%	1.913%
38	18.371	2722	2726	2739	rVB2	406039	1035118	21.42%	1.895%
39	18.748	2786	2790	2796	rBV4	300208	669676	13.86%	1.226%

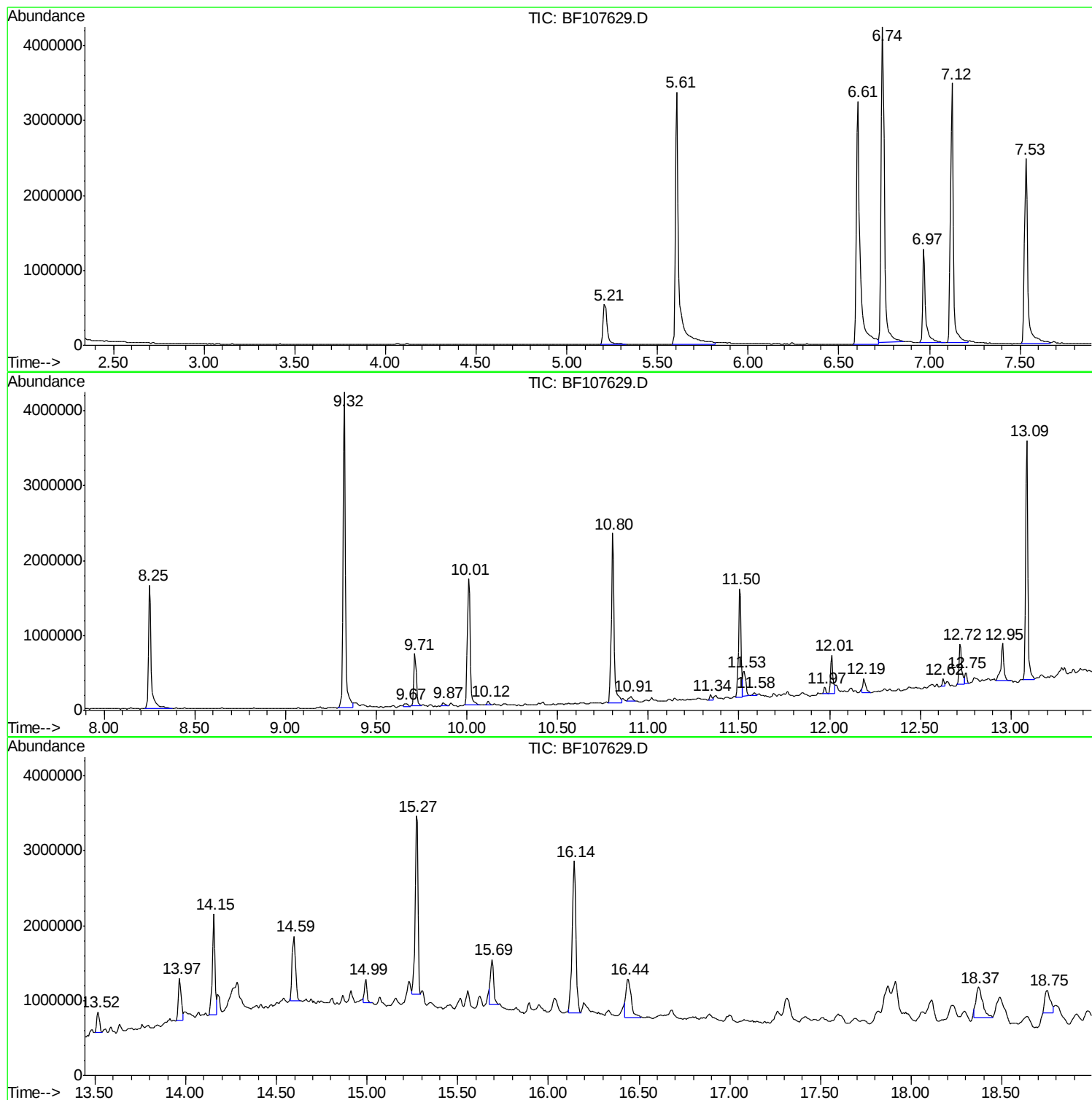
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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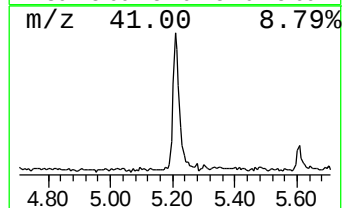
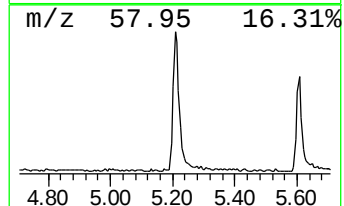
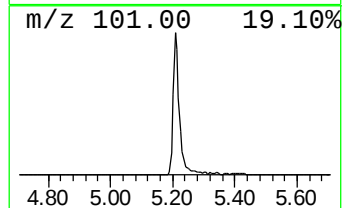
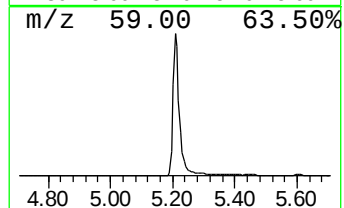
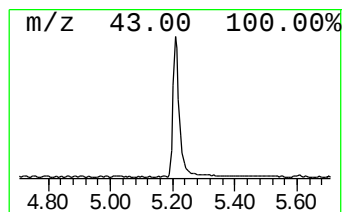
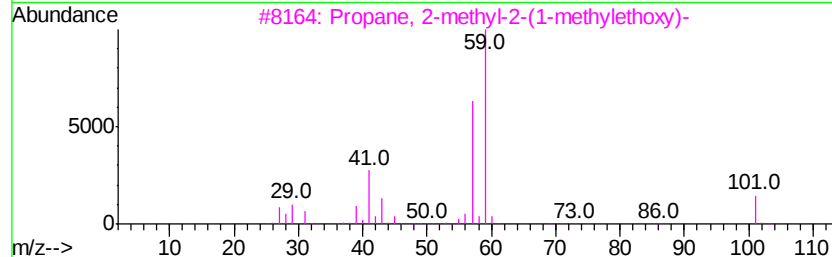
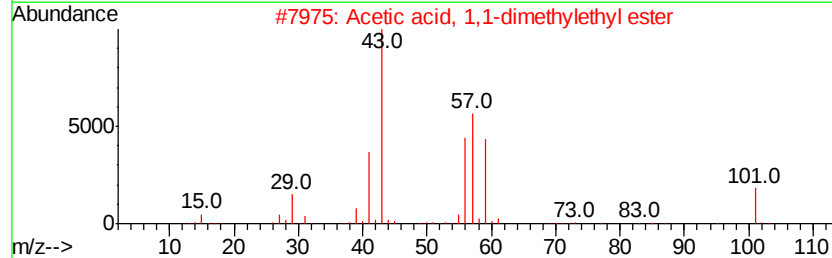
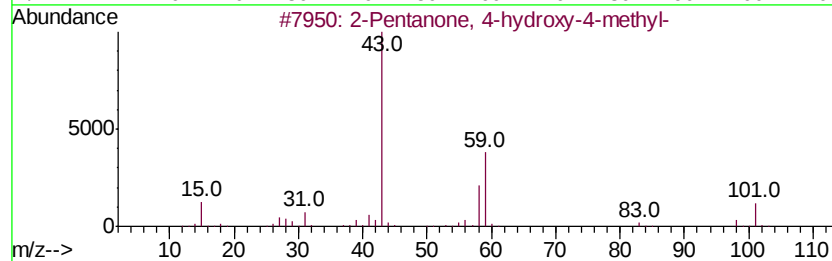
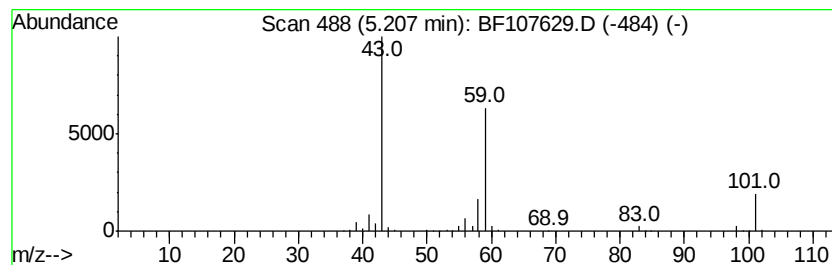
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	10.48 ng	771430	1,4-Dichlorobenzene-d4	6.97

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			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28



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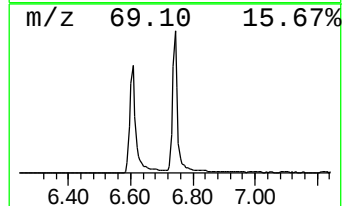
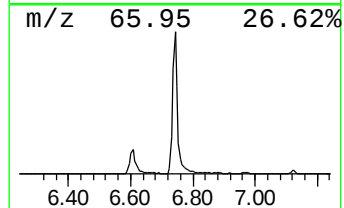
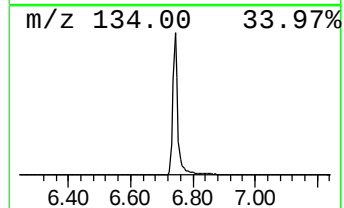
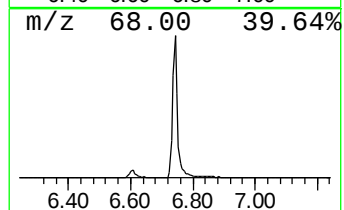
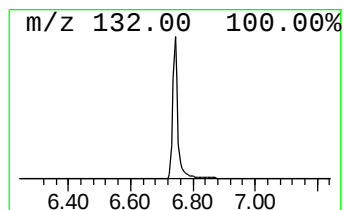
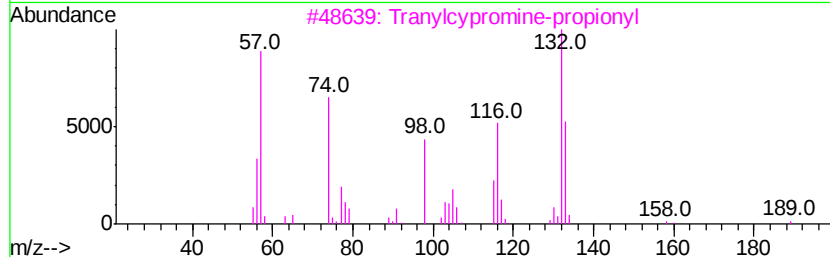
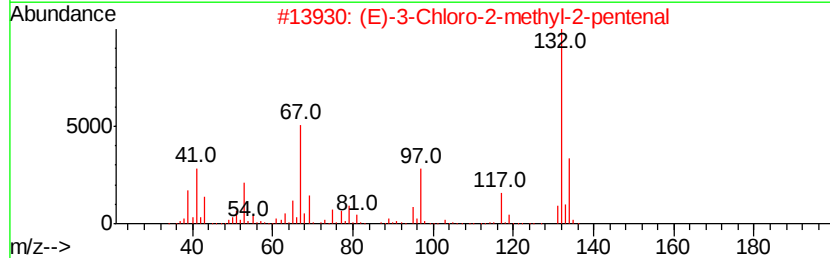
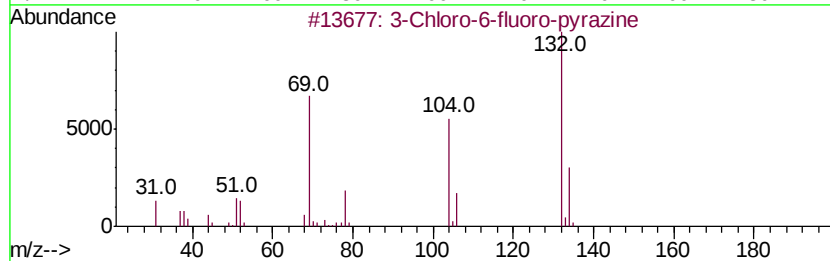
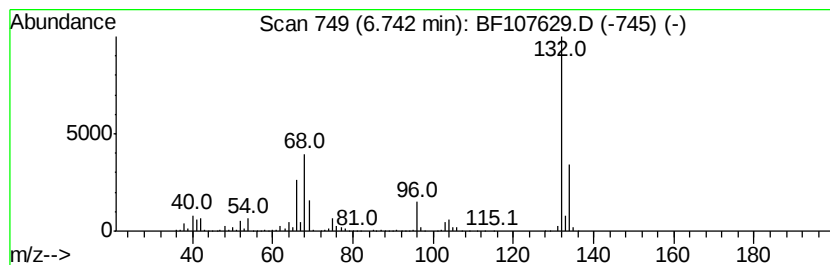
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Peak Number 2 unknown6.74 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	65.66 ng	4831970	1,4-Dichlorobenzene-d4	6.97

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		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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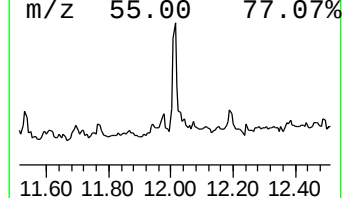
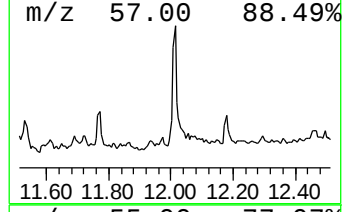
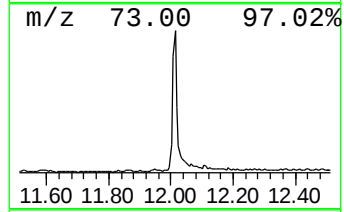
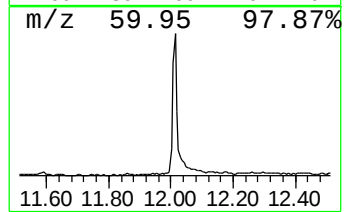
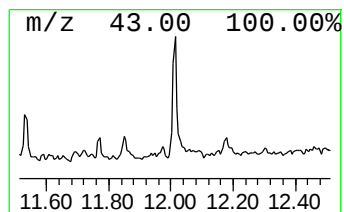
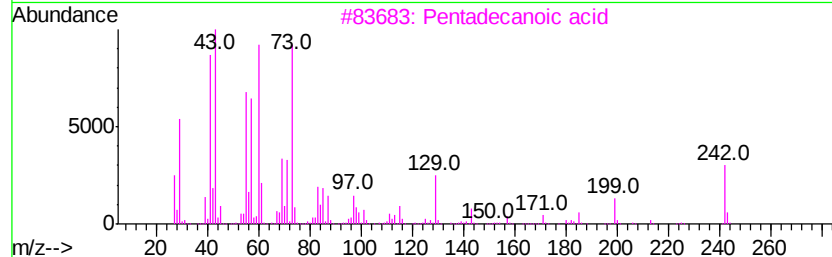
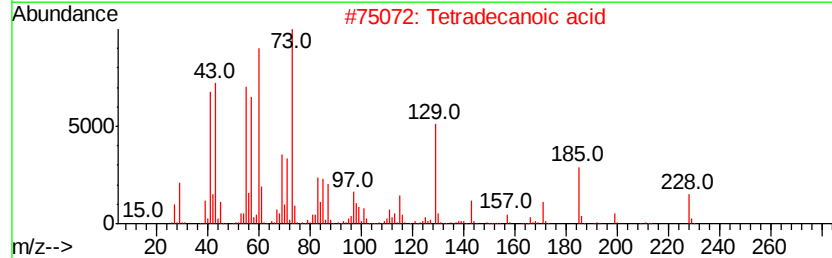
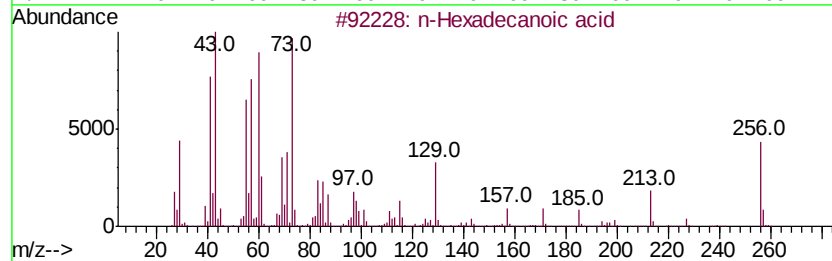
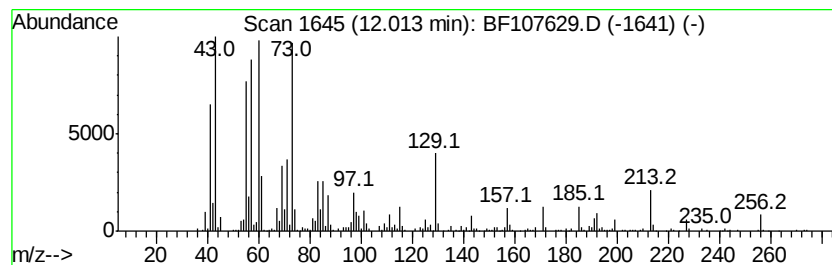
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	7.21 ng	541015	Phenanthrene-d10	11.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	n-Decanoic acid	172	C10H20O2	000334-48-5	60
5	Dodecanoic acid	200	C12H24O2	000143-07-7	58



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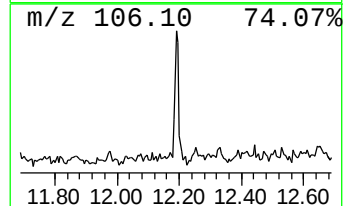
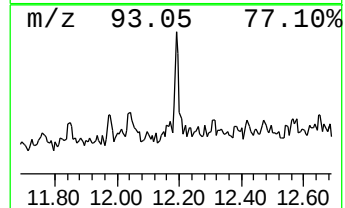
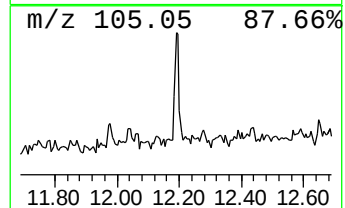
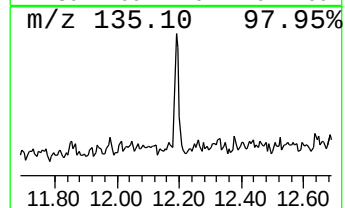
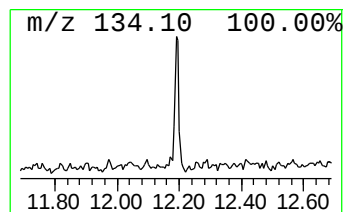
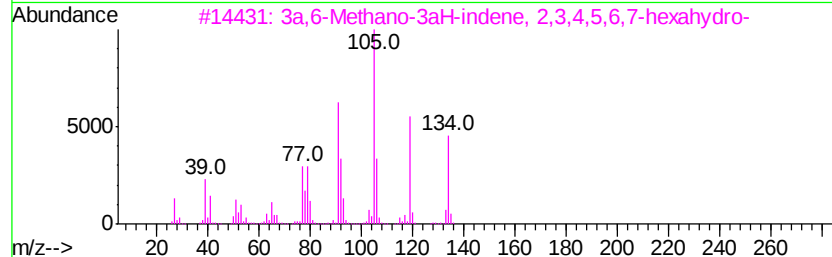
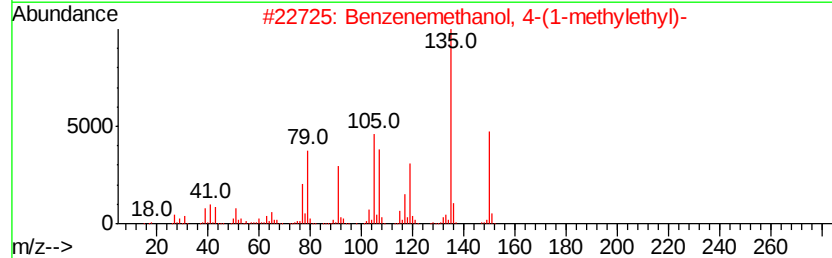
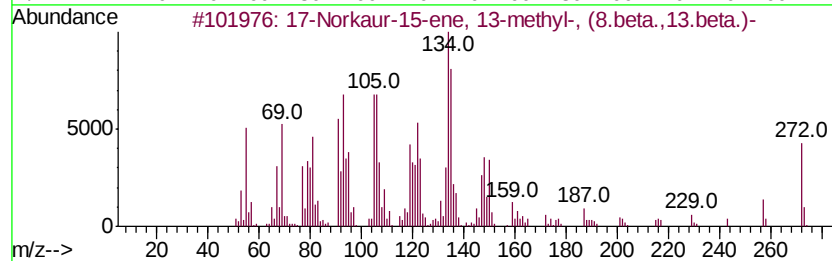
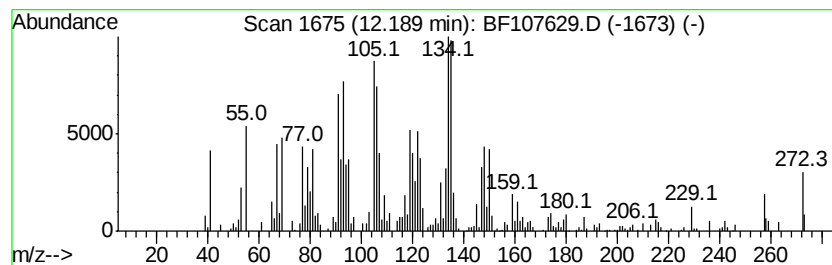
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 17-Norkaur-15-ene, 13-methy... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.19	2.65 ng	199050	Phenanthrene-d10	11.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-Norkaur-15-ene, 13-methyl-, (...	272	C20H32	003564-54-3	96
2		Benzenemethanol, 4-(1-methylethyl)-	150	C10H14O	000536-60-7	30
3		3a,6-Methano-3aH-indene, 2,3,4,5...	134	C10H14	098640-10-9	18
4		Ethanone, 1-(3-methoxyphenyl)-	150	C9H10O2	000586-37-8	18
5		4-Hydroxy-3-methylacetophenone	150	C9H10O2	000876-02-8	18



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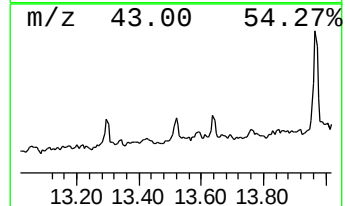
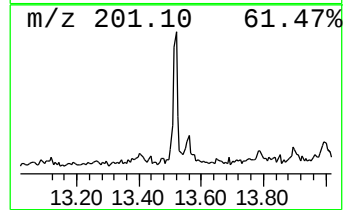
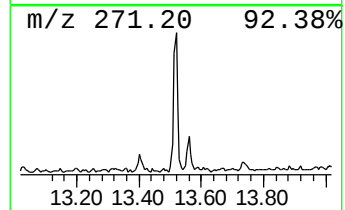
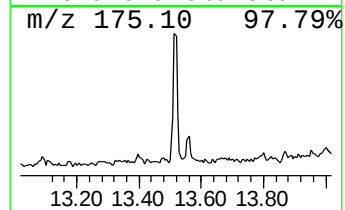
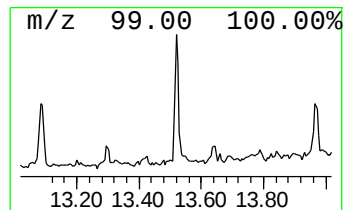
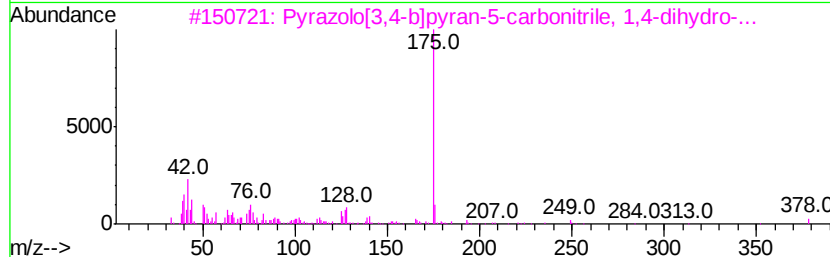
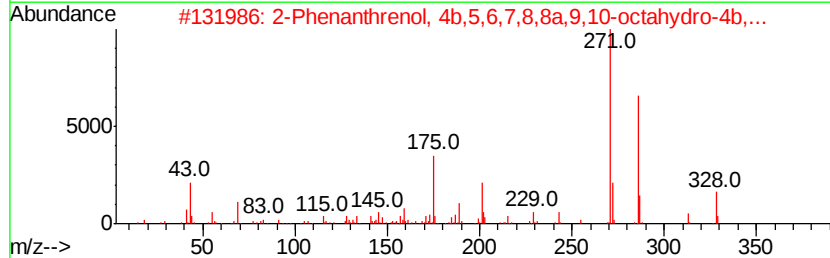
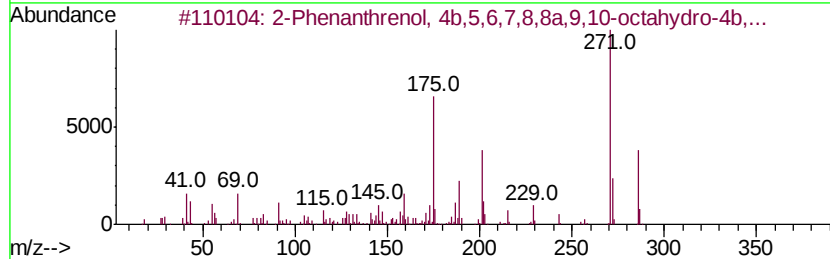
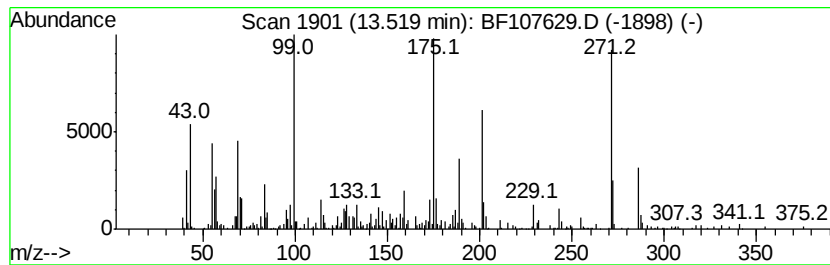
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.52	3.47 ng	253665	Chrysene-d12	14.15		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol,	4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	89
2	2-Phenanthrenol,	4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	43
3	Pvrazolo[3,4-b]p	vrn-5-carbonitr...	378	C14H11IN4O	297162-29-9	10
4	3-Methyl-3-pentyl	methylphosphon...	182	C7H16F02P	1000216-68-4	10
5	4-Methyl-1-pentyl	methylphosphon...	182	C7H16F02P	1000216-68-2	10



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Data File : BF107629.D
Acq On : 24 Jul 2018 19:52
Operator : JU/SJ
Sample : J4098-02
Misc :
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ClientSampleId :
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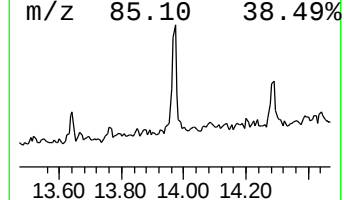
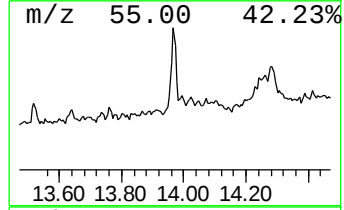
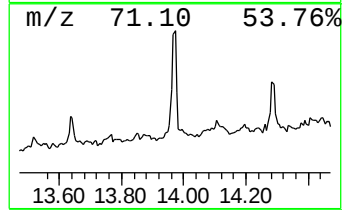
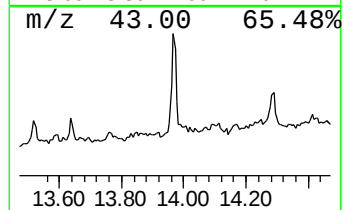
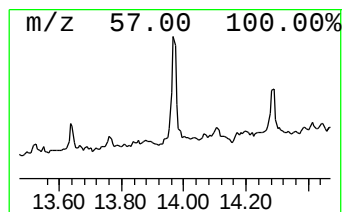
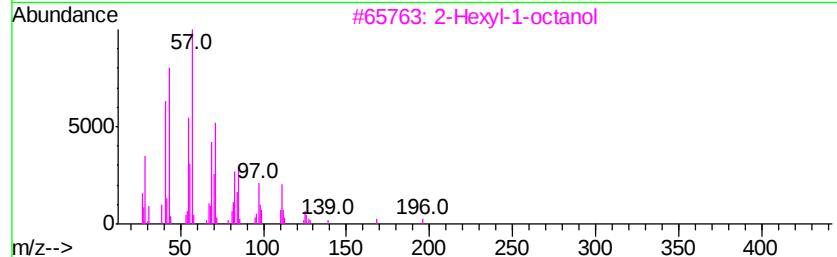
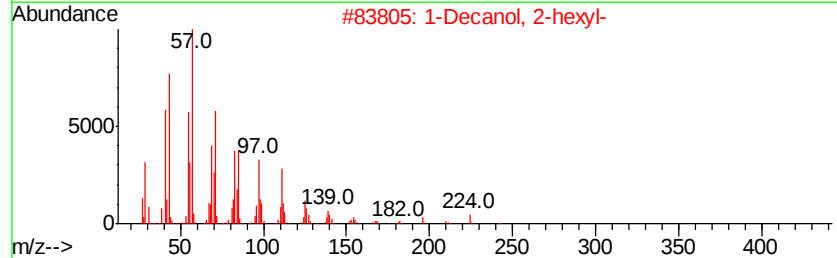
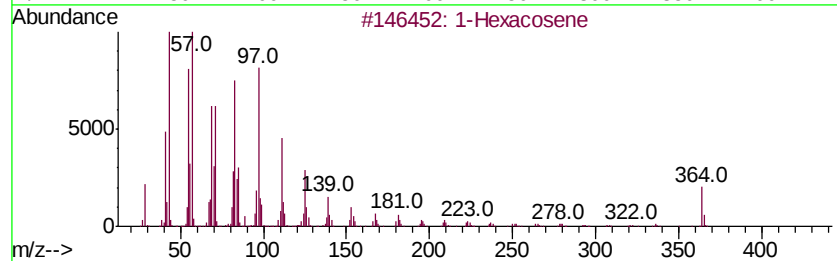
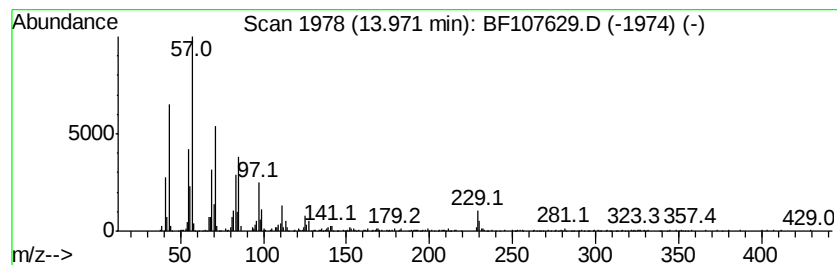
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Hexacosene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	8.15 ng	596172	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexacosene	364	C26H52	018835-33-1	93
2		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	93
3		2-Hexyl-1-octanol	214	C14H30O	019780-79-1	86
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	86
5		Cyclooctacosane	392	C28H56	000297-24-5	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107629.D
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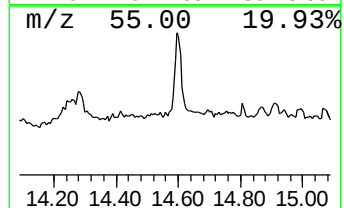
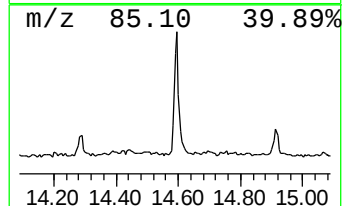
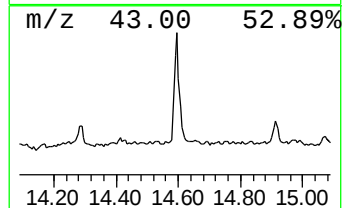
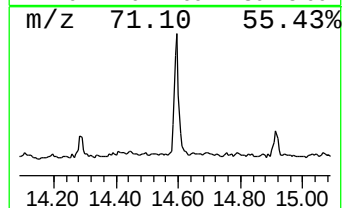
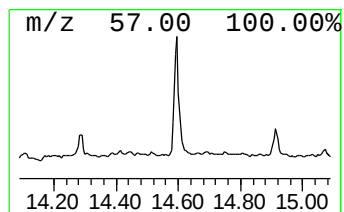
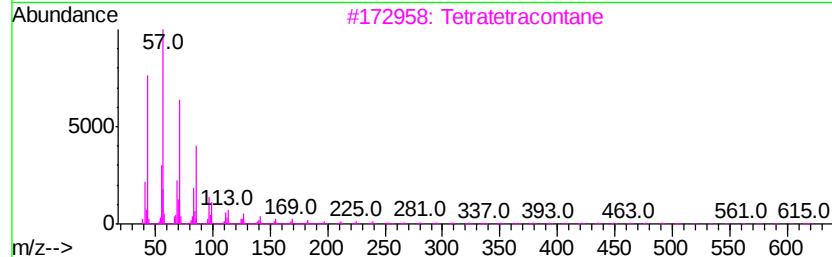
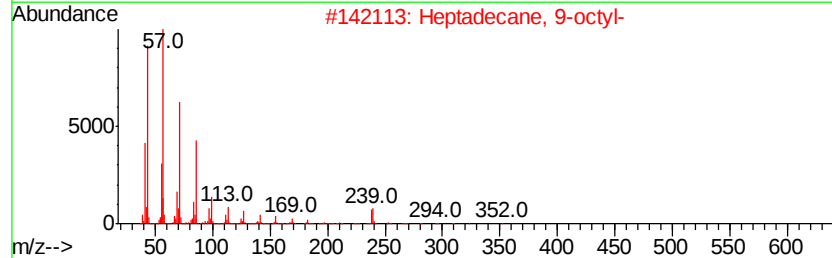
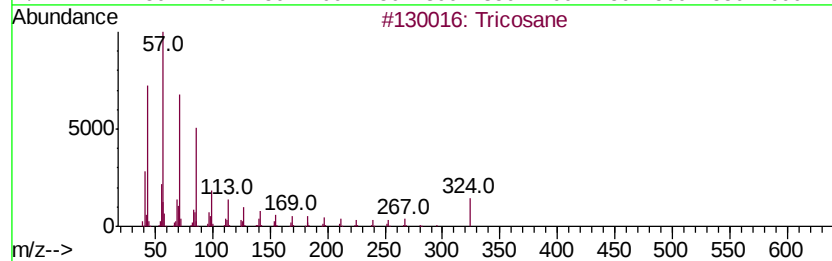
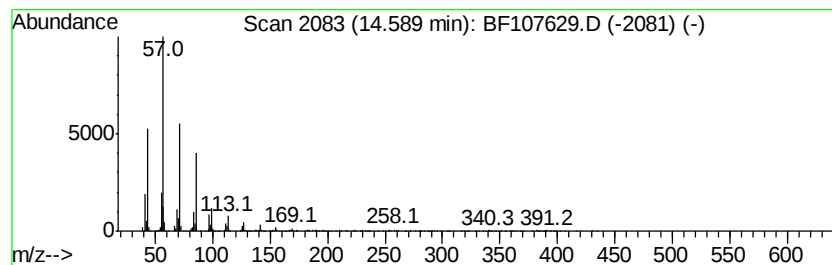
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tricosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	14.08 ng	1030780	Chrysene-d12	14.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane		324	C23H48	000638-67-5	94
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	93
3	Tetratetracontane		619	C44H90	007098-22-8	91
4	Tritetracontane		605	C43H88	007098-21-7	91
5	Eicosane, 10-methyl-		296	C21H44	054833-23-7	91



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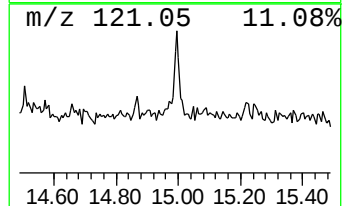
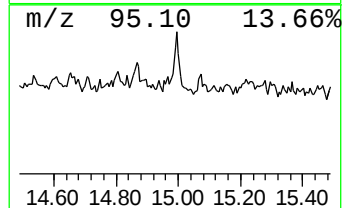
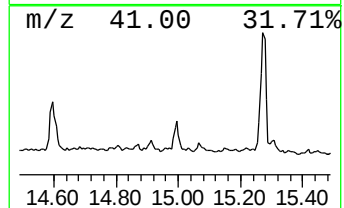
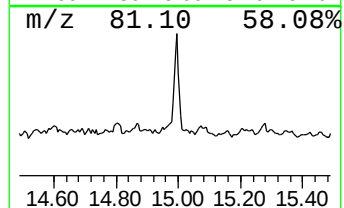
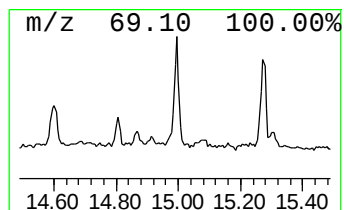
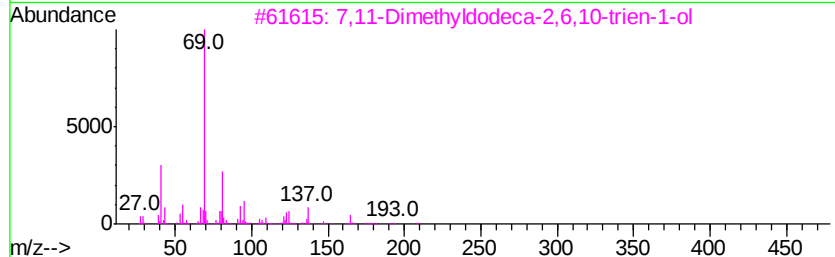
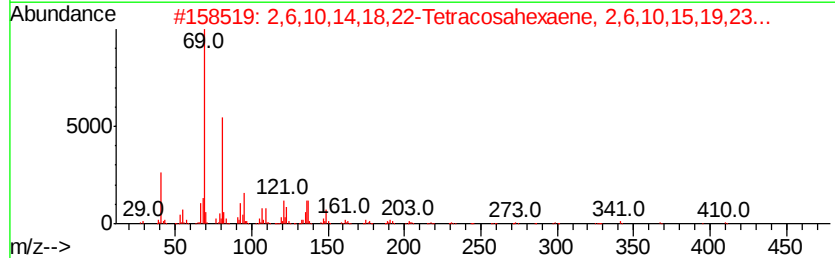
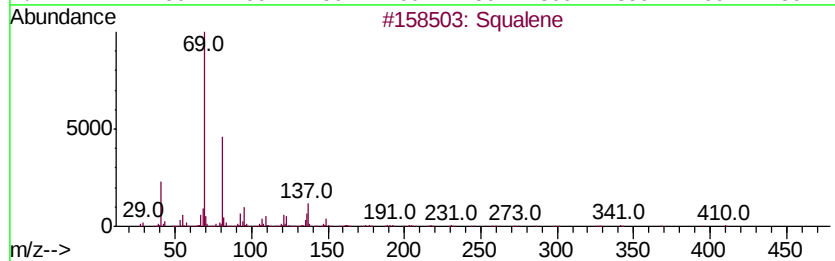
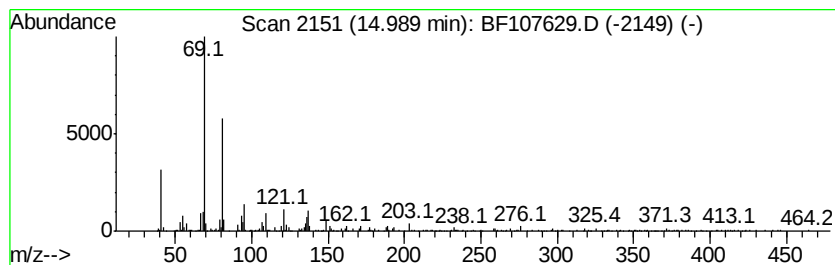
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Squalene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	8.32 ng	337744	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	98
2		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	93
3		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
4		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	64
5		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	000106-28-5	56



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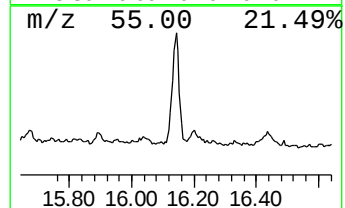
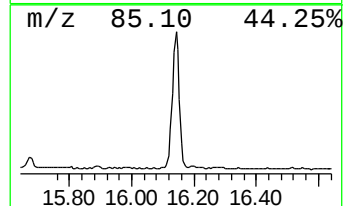
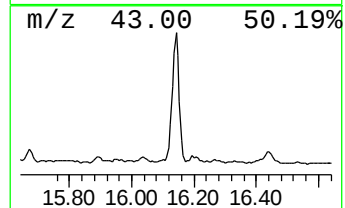
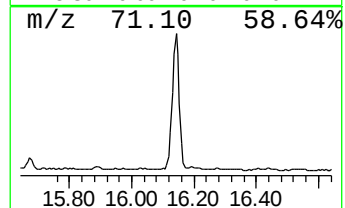
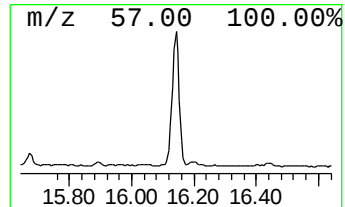
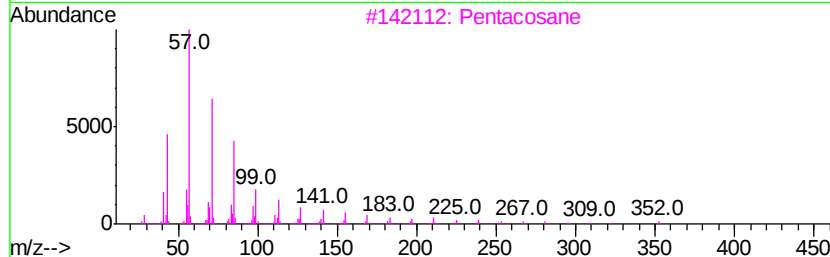
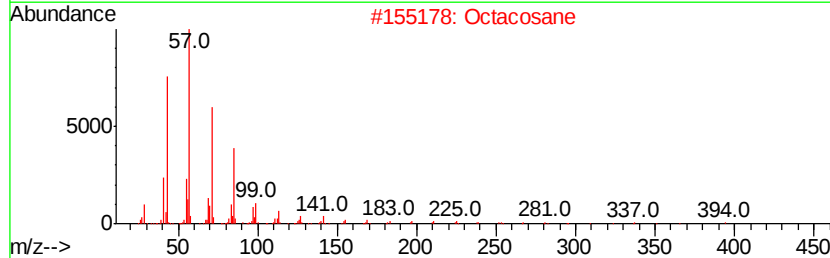
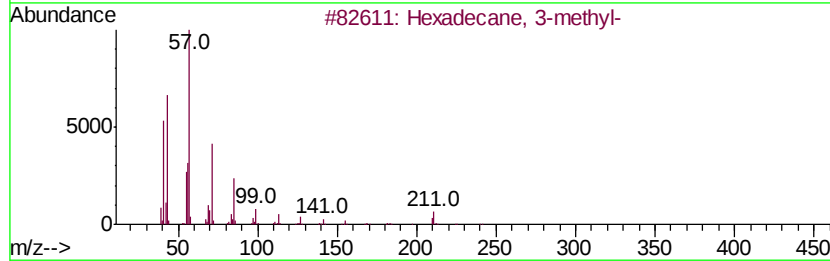
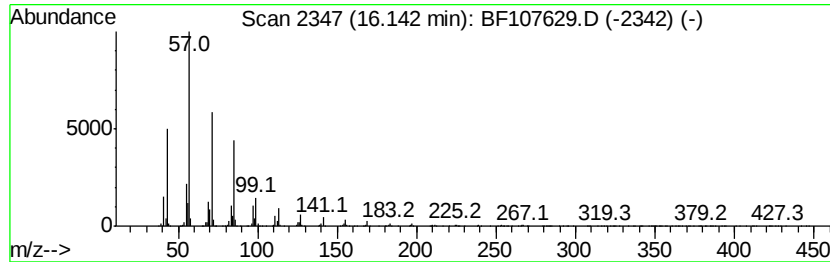
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Hexadecane, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	76.18 ng	3092670	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Heptacosane	380	C27H56	000593-49-7	91



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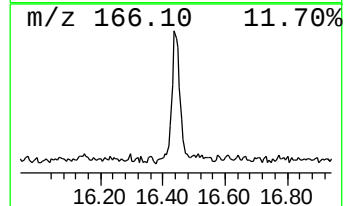
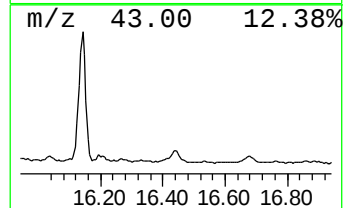
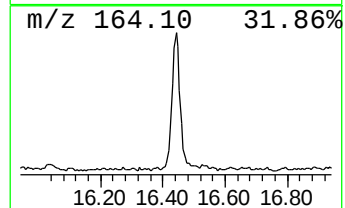
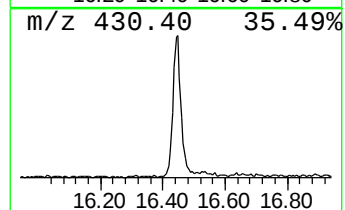
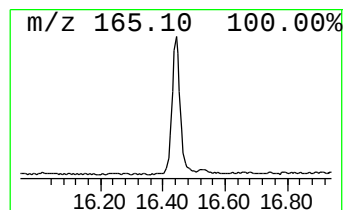
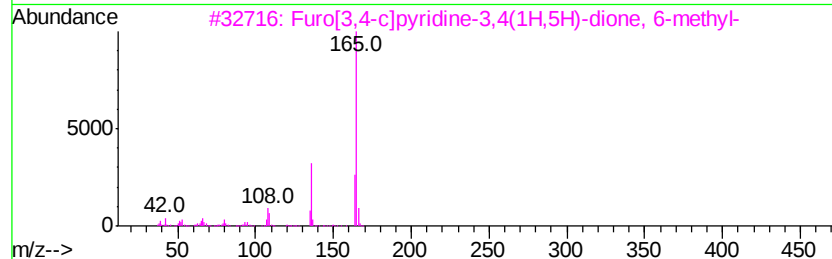
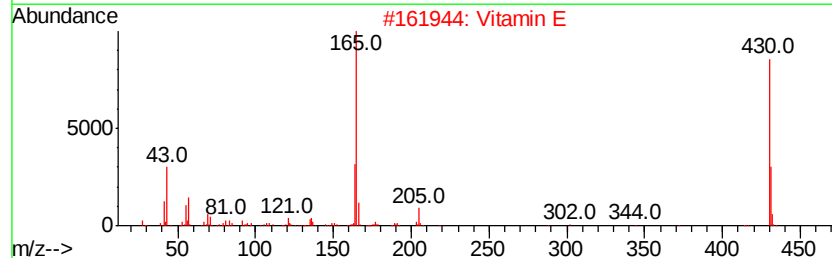
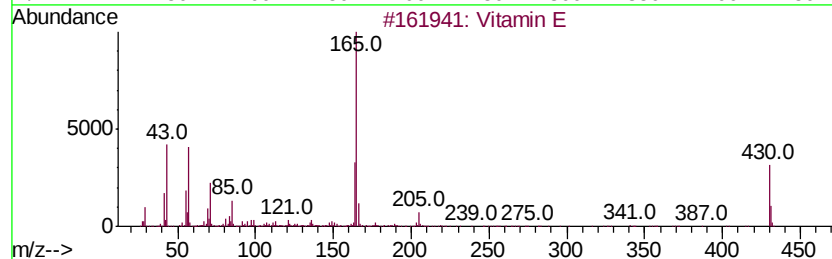
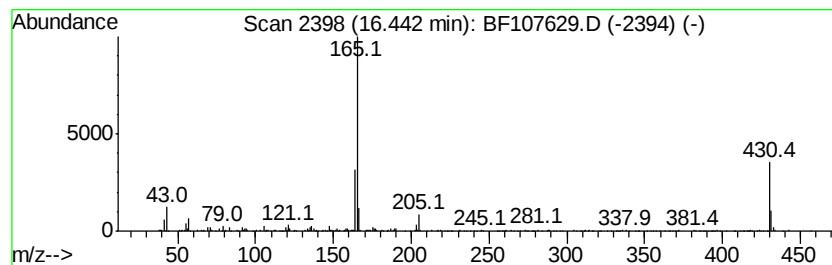
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Vitamin E Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	25.74 ng	1044890	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vitamin E	430	C29H50O2	000059-02-9	95
2		Vitamin E	430	C29H50O2	010191-41-0	93
3		Furo[3,4-c]pyridine-3,4(1H,5H)-d...	165	C8H7N03	007472-18-6	58
4		Piperazine, 1-(3,4-dimethoxybenz...	430	C26H26N2O4	333756-87-9	50
5		Isoxaben	332	C18H24N2O4	082558-50-7	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
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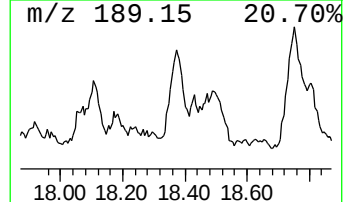
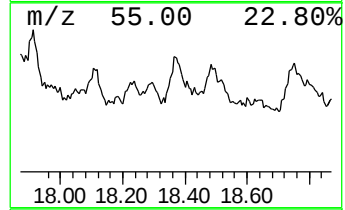
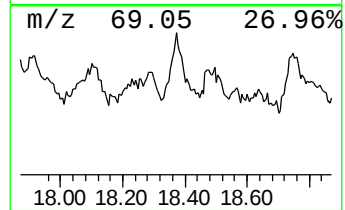
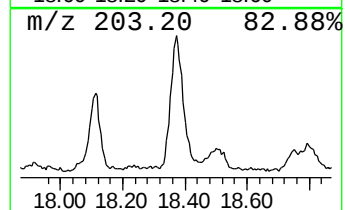
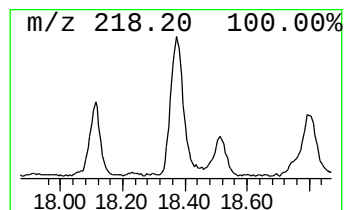
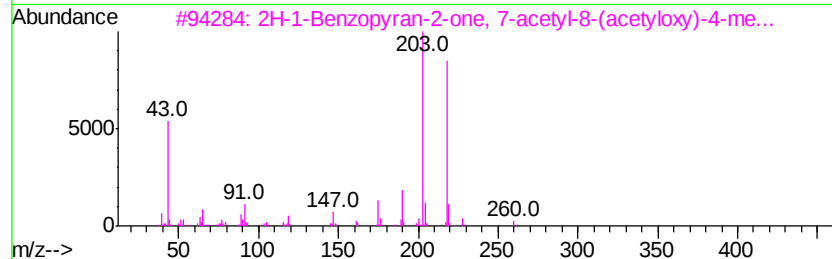
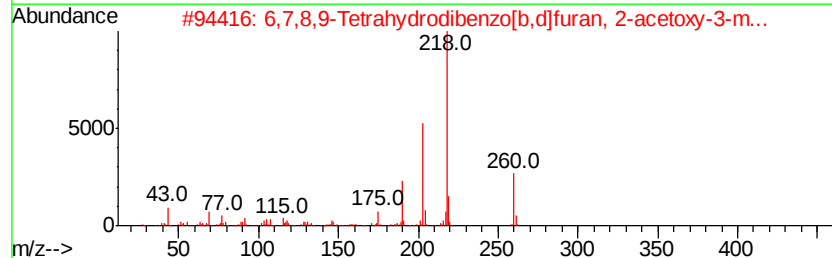
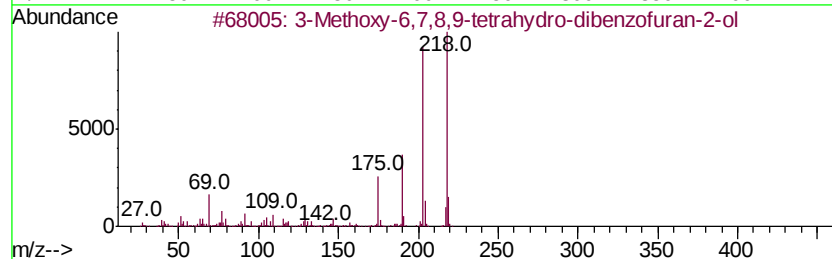
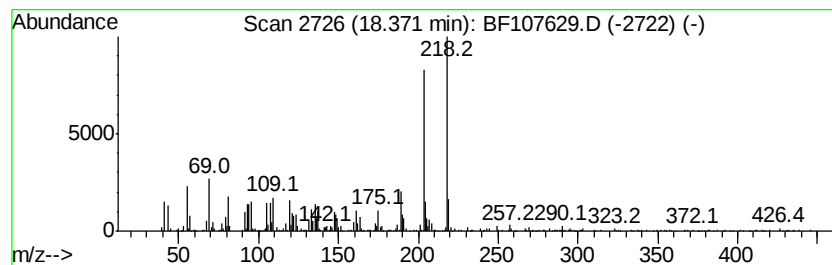
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 3-Methoxy-6,7,8,9-tetrahydr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	25.50 ng	1035120	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Methoxy-6,7,8,9-tetrahydro-dib...	218	C13H14O3	1000186-57-9	68
2		6,7,8,9-Tetrahydrodibenzo[b,d]fu...	260	C15H16O4	1000196-34-3	64
3		2H-1-Benzopyran-2-one, 7-acetyl-...	260	C14H12O5	129812-32-4	47
4		5(1H)-Azulenone, 2,4,6,7,8,8a-he...	218	C15H22O	006754-66-1	46
5		Digitoxinigenin-4-en	372	C23H32O4	001778-88-7	46



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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.74	6.74	65.7	ng	4831970	1	6.97	1471760	20.0
n-Hexadecanoic acid	12.01	7.2	ng	541015	4	11.50	1500270	20.0
17-Norkaur-15-ene...	12.19	2.6	ng	199050	4	11.50	1500270	20.0
2-Phenanthrenol, ...	13.52	3.5	ng	253665	5	14.15	1463760	20.0
1-Hexacosene	13.97	8.2	ng	596172	5	14.15	1463760	20.0
Tricosane	14.59	14.1	ng	1030780	5	14.15	1463760	20.0
Squalene	14.99	8.3	ng	337744	6	15.69	811972	20.0
Hexadecane, 3-met...	16.14	76.2	ng	3092670	6	15.69	811972	20.0
Vitamin E	16.44	25.7	ng	1044890	6	15.69	811972	20.0
3-Methoxy-6,7,8,9...	18.37	25.5	ng	1035120	6	15.69	811972	20.0
5-Bromo-4-oxo-4,5...	18.75	16.5	ng	669676	6	15.69	811972	20.0