

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121518\
 Data File : BF111459.D
 Acq On : 15 Dec 2018 10:55
 Operator : JU/SJ
 Sample : J6208-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121418.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	3.240	193	196	218	rBV	81546	187986	2.16%	0.223%
2	5.004	493	496	502	rVB	166616	223120	2.57%	0.265%
3	5.422	562	567	571	rBV	4025570	4236070	48.72%	5.030%
4	6.439	735	740	757	rBV	4069645	4878047	56.10%	5.792%
5	6.569	757	762	765	rBV	4802752	4438802	51.05%	5.271%
6	6.792	796	800	804	rBV	1159972	922588	10.61%	1.095%
7	6.951	823	827	830	rBV	3821617	3215886	36.98%	3.819%
8	7.357	890	896	899	rBV	2921869	2860316	32.89%	3.396%
9	8.075	1014	1018	1021	rBV	1494028	1168656	13.44%	1.388%
10	9.151	1196	1201	1204	rBV	4628767	3994655	45.94%	4.743%
11	9.545	1264	1268	1271	rBV	488438	415115	4.77%	0.493%
12	9.827	1310	1316	1320	rBV	1504678	1335987	15.36%	1.586%
13	10.045	1347	1353	1360	rBV2	186350	233508	2.69%	0.277%
14	10.621	1446	1451	1454	rBV	2766822	2741952	31.53%	3.256%
15	10.986	1509	1513	1521	rVB	220585	265178	3.05%	0.315%
16	11.316	1565	1569	1571	rBV	1371224	1327035	15.26%	1.576%
17	11.333	1571	1572	1576	rVV	320460	244933	2.82%	0.291%
18	11.527	1592	1605	1606	rBV	507956	1353432	15.56%	1.607%
19	11.704	1633	1635	1643	rBV5	44156	88299	1.02%	0.105%
20	11.780	1643	1648	1649	rBV4	103573	138422	1.59%	0.164%
21	11.874	1651	1664	1670	rVV	4670648	8695445	100.00%	10.325%
22	12.521	1771	1774	1777	rBV	632539	669071	7.69%	0.794%
23	12.562	1778	1781	1788	rVV2	687965	999107	11.49%	1.186%
24	12.645	1788	1795	1802	rVV	3294126	4930057	56.70%	5.854%
25	12.727	1803	1809	1811	rVV2	667421	1364305	15.69%	1.620%
26	12.751	1811	1813	1817	rVB	674802	614296	7.06%	0.729%
27	12.857	1823	1831	1835	rBV	1687781	3837388	44.13%	4.557%
28	12.904	1835	1839	1842	rVB	3397625	3370291	38.76%	4.002%
29	13.021	1853	1859	1864	rVB3	166928	356378	4.10%	0.423%
30	13.133	1875	1878	1882	rVB3	151119	202452	2.33%	0.240%
31	13.351	1912	1915	1922	rBV4	122117	228949	2.63%	0.272%
32	13.451	1929	1932	1935	rBV	91874	99669	1.15%	0.118%
33	13.645	1957	1965	1970	rBV	2828903	4740340	54.52%	5.629%
34	13.804	1988	1992	2000	rVB3	485553	721713	8.30%	0.857%

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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-BF121418.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.951	2012	2017	2020	rBV2	1125174	1257937	14.47%	1.494%
36	13.974	2020	2021	2024	rVB	204954	113859	1.31%	0.135%
37	14.239	2058	2066	2071	rBV	2892249	4474767	51.46%	5.313%
38	14.427	2095	2098	2105	rBV5	94555	172460	1.98%	0.205%
39	14.545	2115	2118	2125	rVB2	87658	132034	1.52%	0.157%
40	14.739	2144	2151	2157	rBV	2385702	3943295	45.35%	4.682%
41	14.974	2187	2191	2199	rVB	201061	373382	4.29%	0.443%
42	15.056	2201	2205	2215	rBV2	233111	600853	6.91%	0.713%
43	15.292	2237	2245	2249	rBV	1618218	2754828	31.68%	3.271%
44	15.327	2249	2251	2257	rVV	192338	200691	2.31%	0.238%
45	15.392	2257	2262	2265	rVV	712459	876582	10.08%	1.041%
46	15.427	2265	2268	2273	rVB2	209257	293301	3.37%	0.348%
47	15.851	2336	2340	2342	rBV	157470	221397	2.55%	0.263%
48	15.898	2342	2348	2363	rVB	729272	1763598	20.28%	2.094%
49	16.621	2464	2471	2480	rBV	423879	955616	10.99%	1.135%
50	17.227	2570	2574	2581	rVB2	67550	135102	1.55%	0.160%
51	17.462	2607	2614	2626	rVB3	219124	675891	7.77%	0.803%
52	17.592	2632	2636	2647	rVB3	74004	171445	1.97%	0.204%

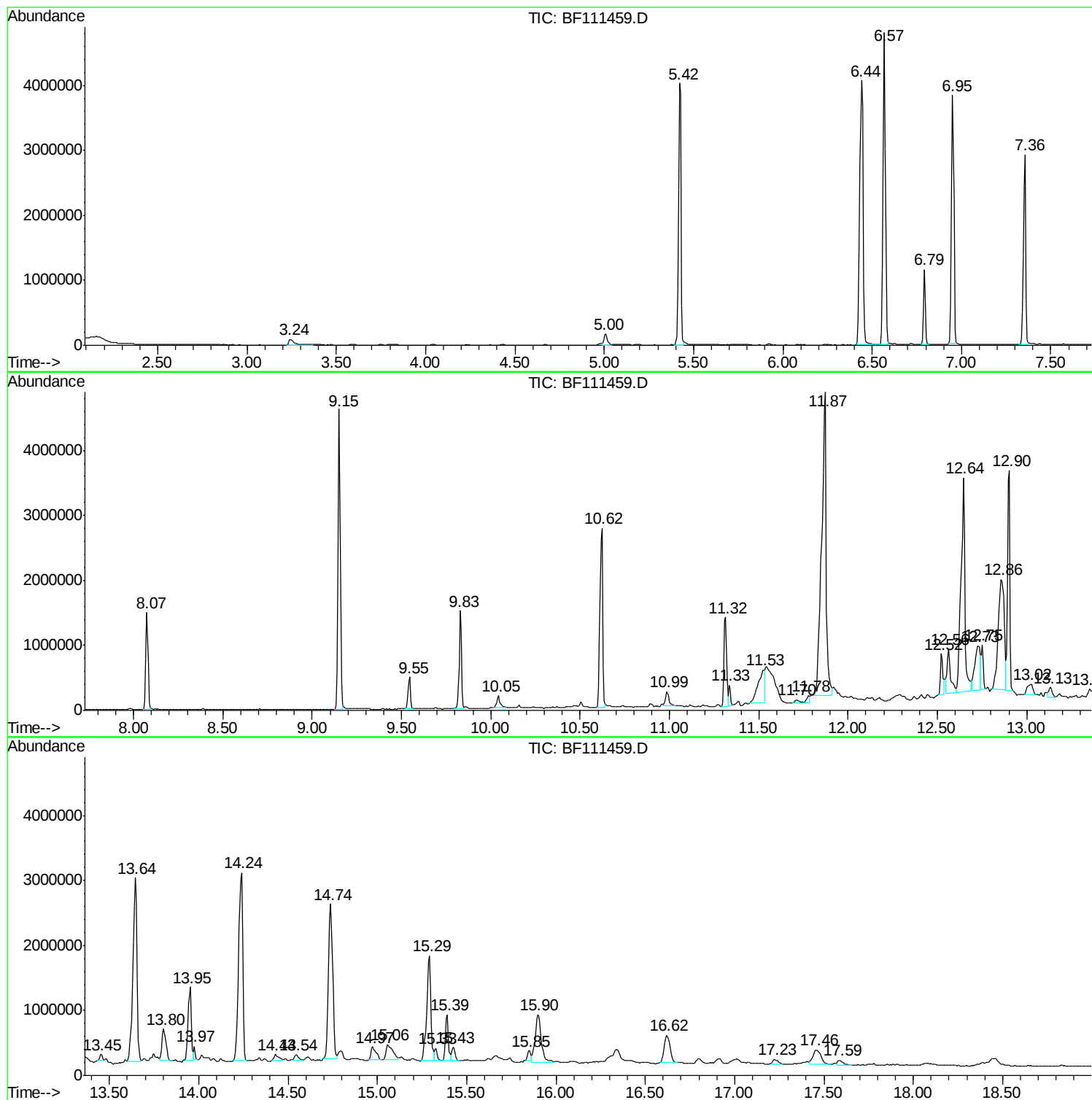
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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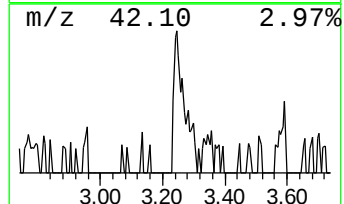
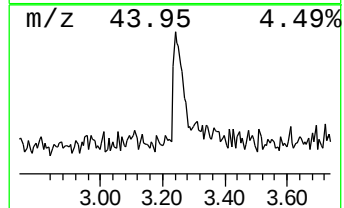
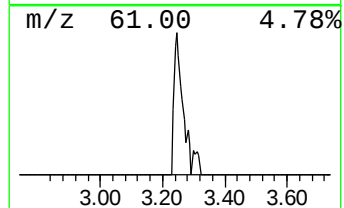
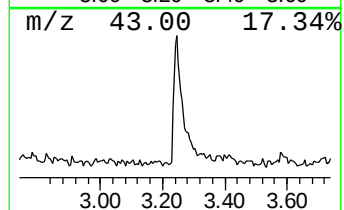
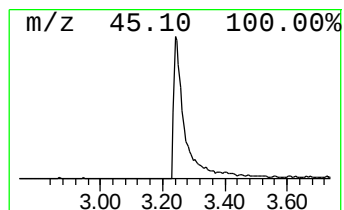
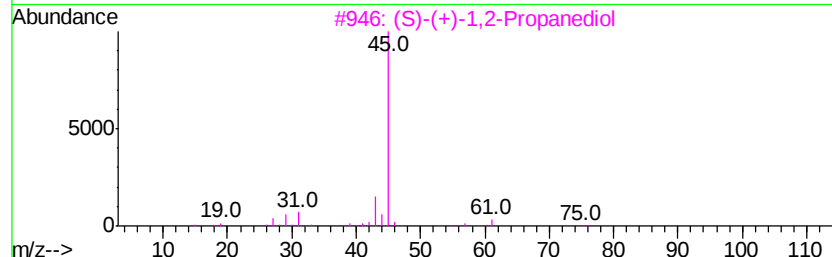
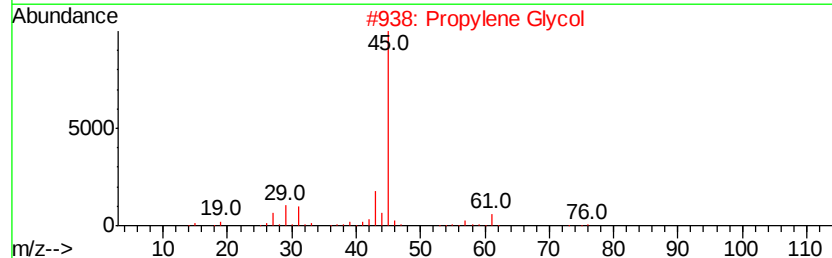
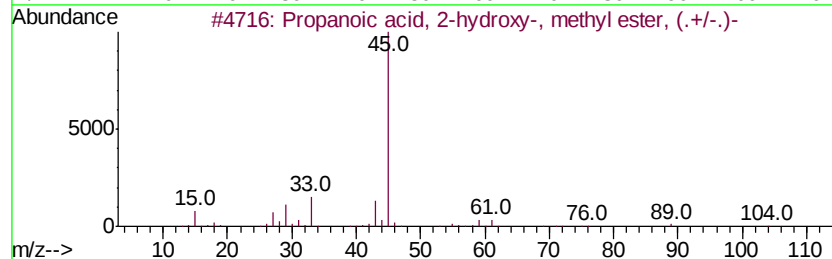
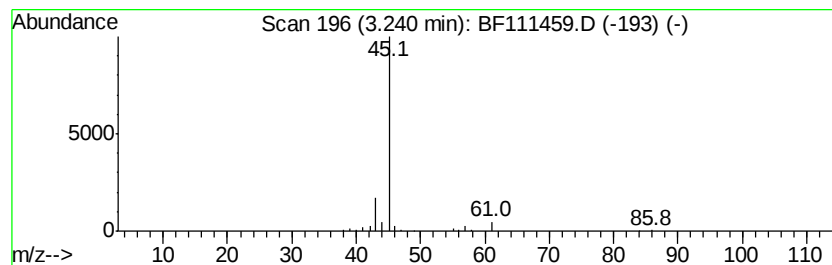
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.24	4.08 ng	187986	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	002155-30-8	43
2			Propylene Glycol	76	C3H8O2	000057-55-6	43
3			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9



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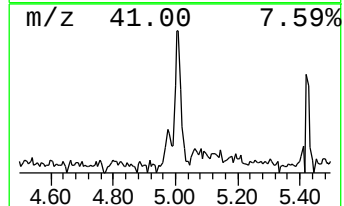
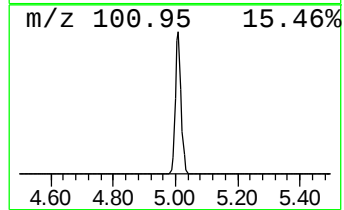
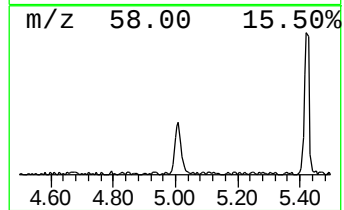
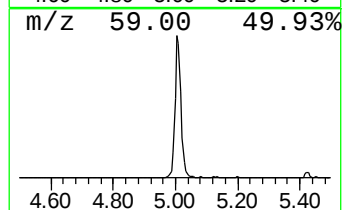
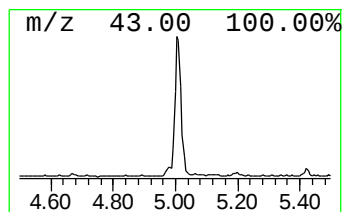
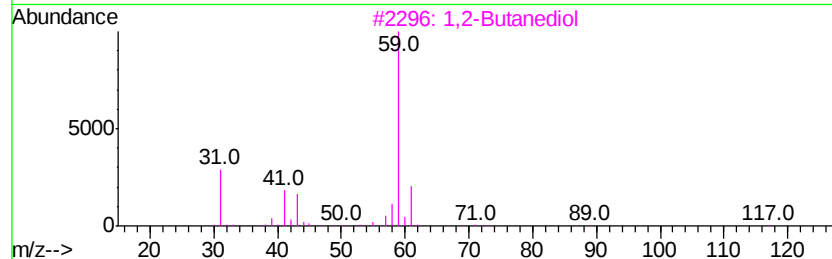
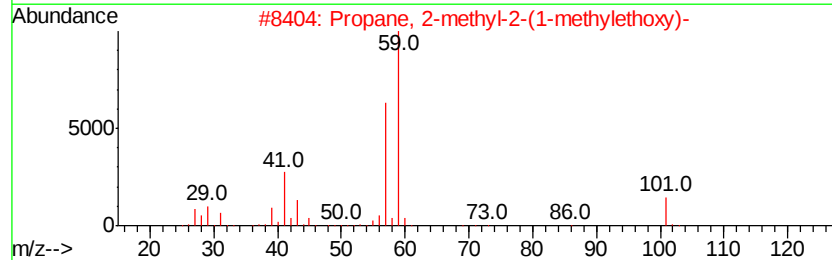
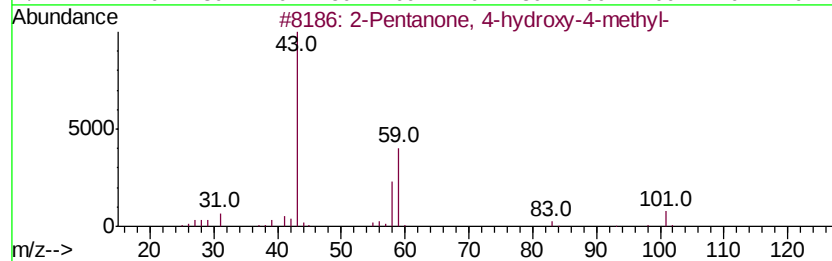
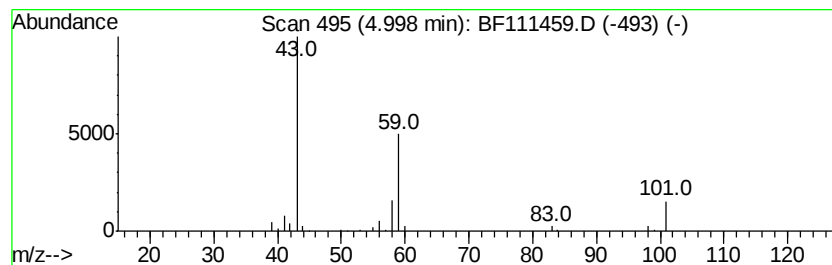
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	4.84 ng	223120	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	42
3		1,2-Butanediol	90	C4H10O2	000584-03-2	28
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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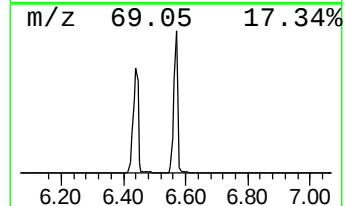
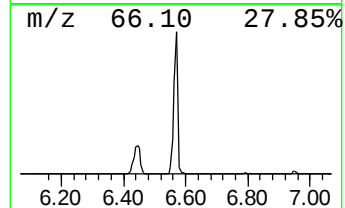
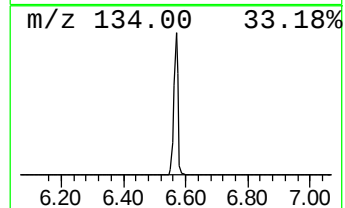
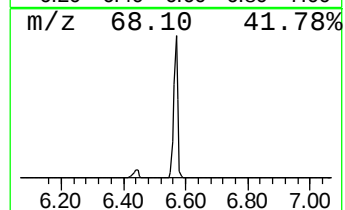
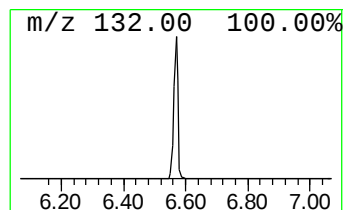
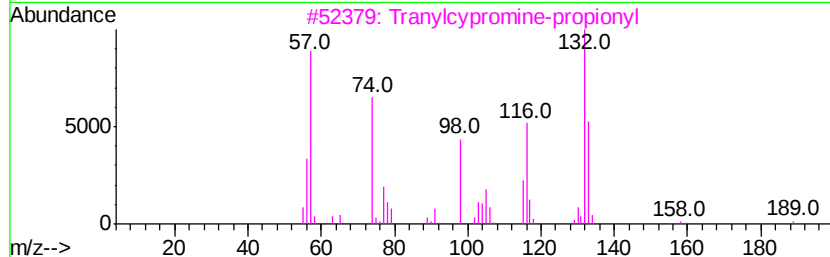
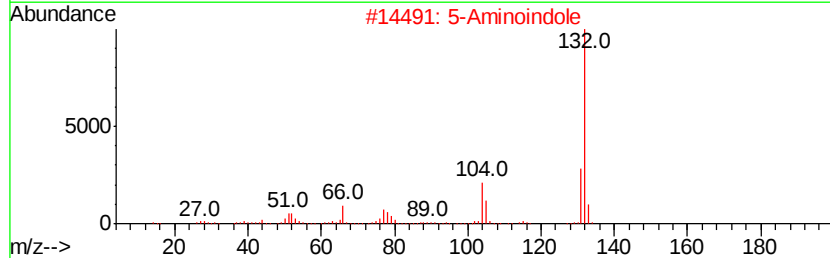
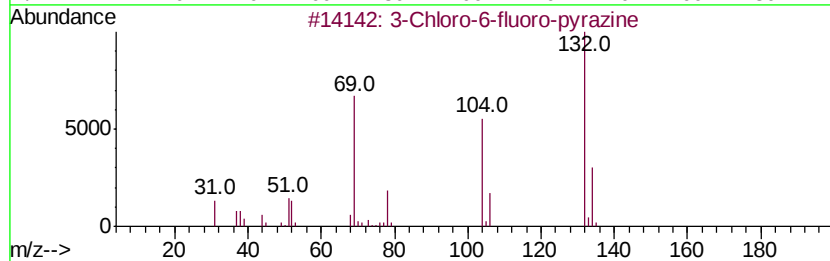
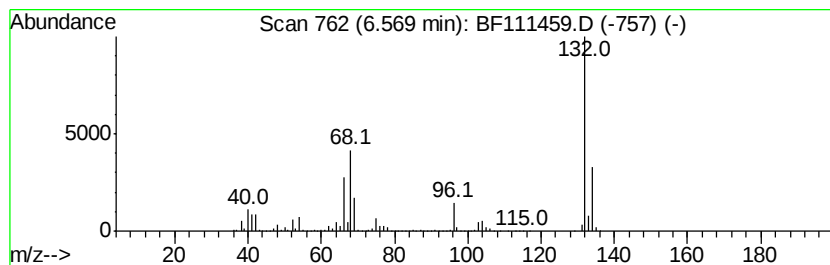
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	96.23 ng	4438800	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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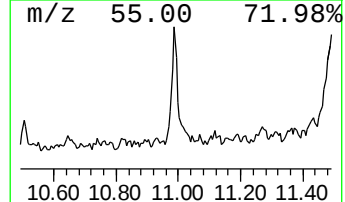
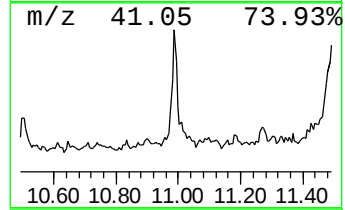
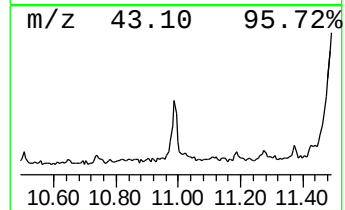
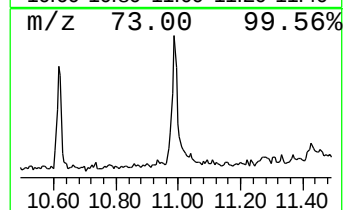
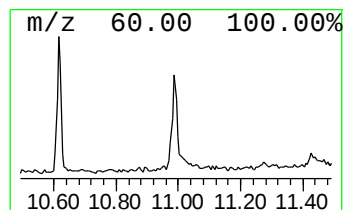
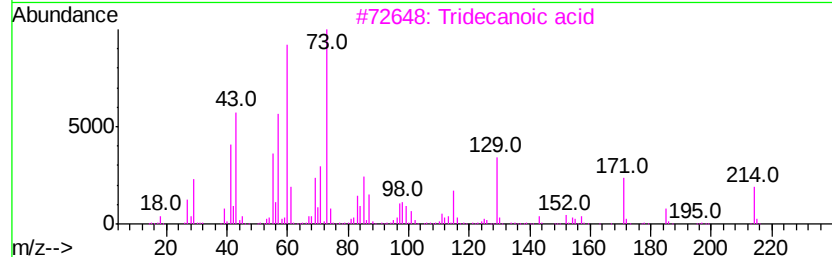
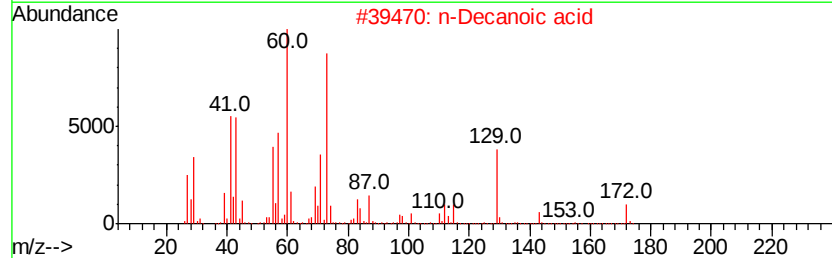
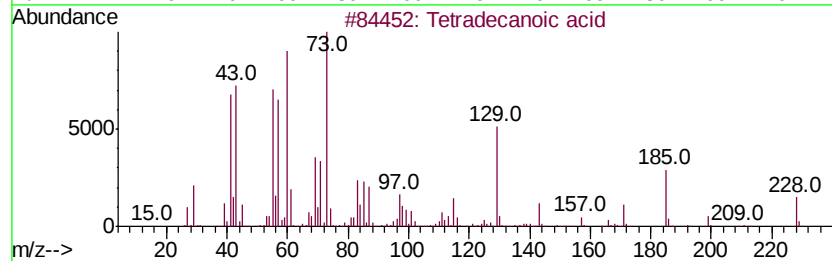
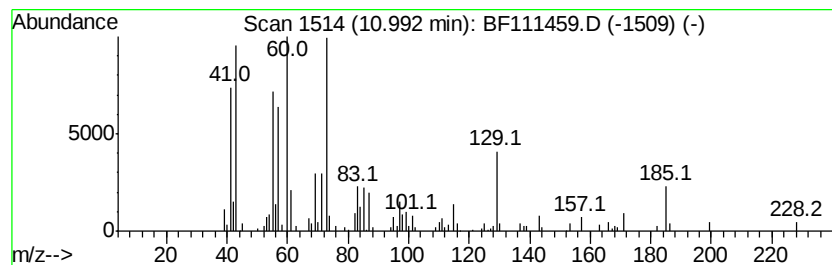
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	4.00 ng	265178	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
2		n-Decanoic acid	172	C10H20O2	000334-48-5	80
3		Tridecanoic acid	214	C13H26O2	000638-53-9	72
4		Dodecanoic acid	200	C12H24O2	000143-07-7	72
5		.beta.-D-Mannofuranoside, 1-O-(1...	332	C17H32O6	1000155-29-9	11



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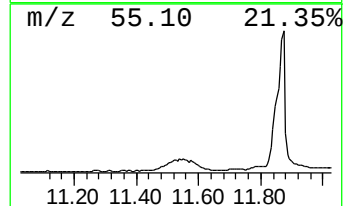
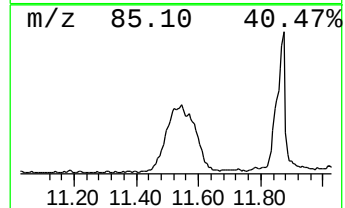
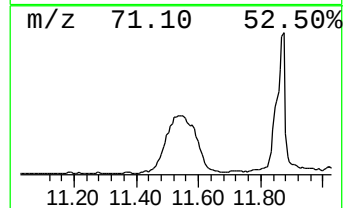
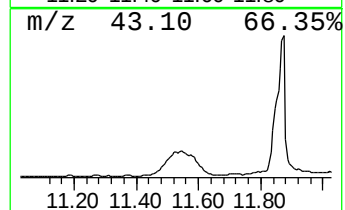
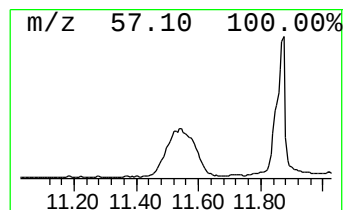
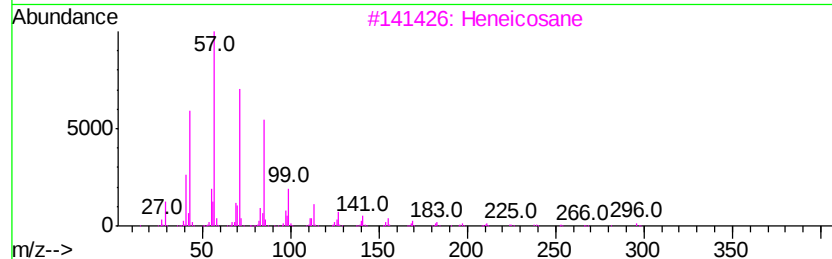
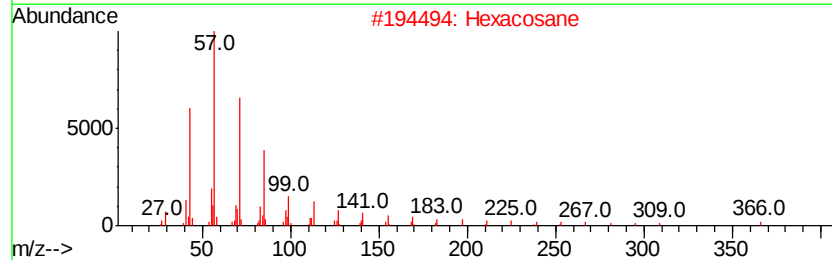
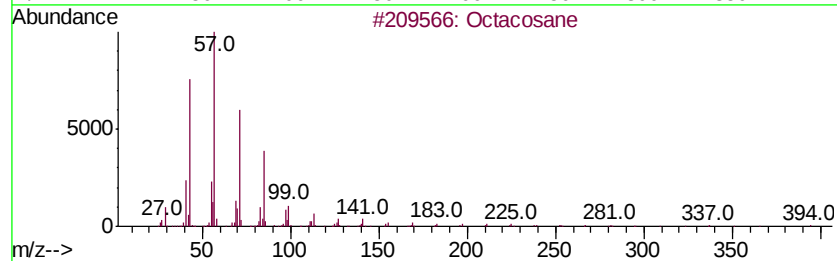
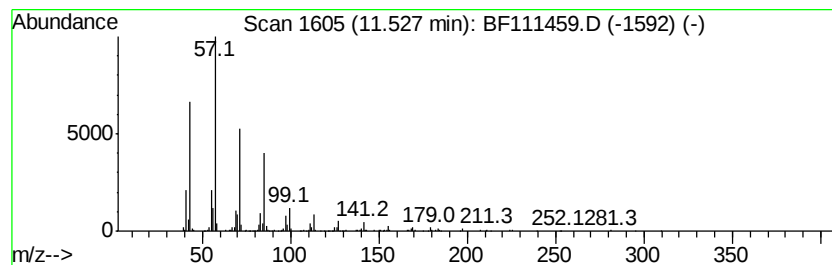
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	20.40 ng	1353430	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Pentacosane	352	C25H52	000629-99-2	90
5		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111459.D
Acq On : 15 Dec 2018 10:55
Operator : JU/SJ
Sample : J6208-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

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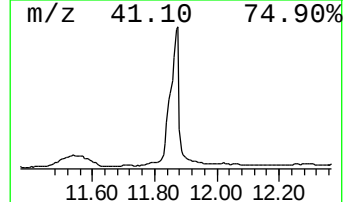
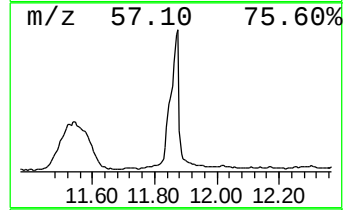
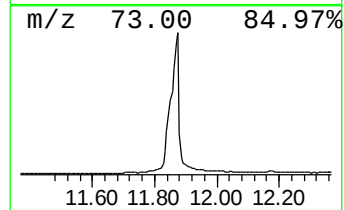
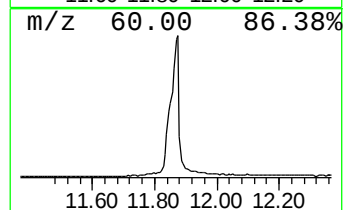
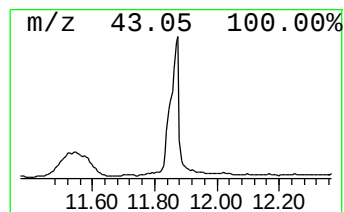
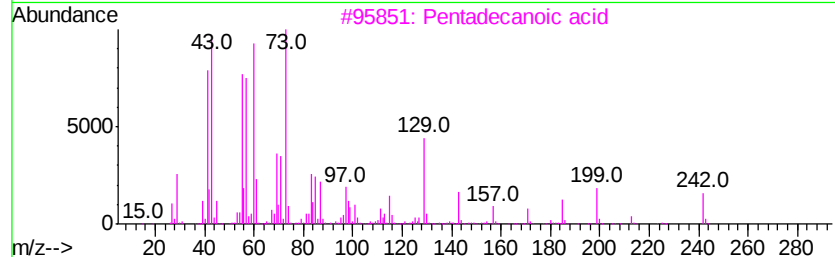
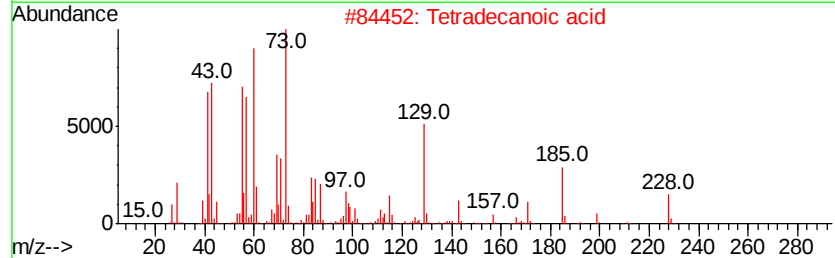
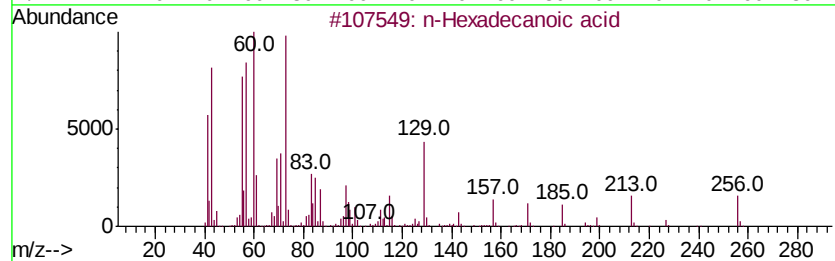
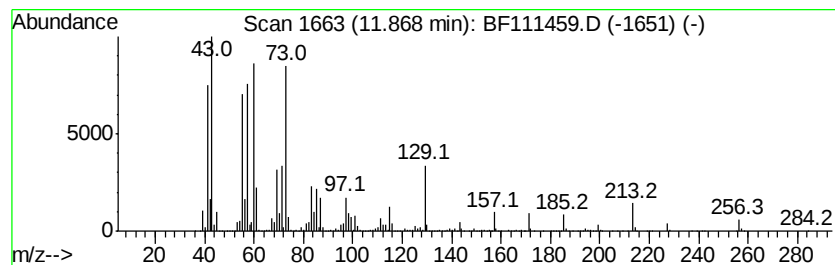
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	131.05 ng	8695450	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tridecanoic acid	214	C13H26O2	000638-53-9	72
5		n-Decanoic acid	172	C10H20O2	000334-48-5	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
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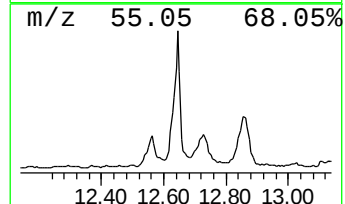
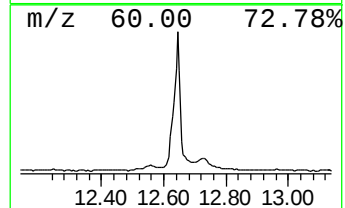
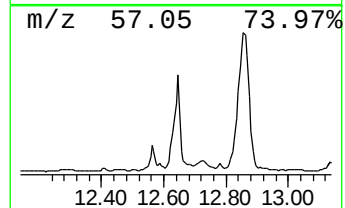
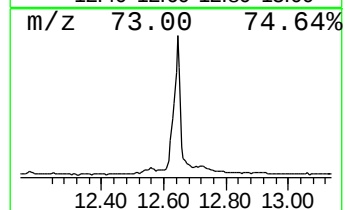
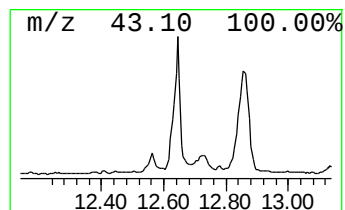
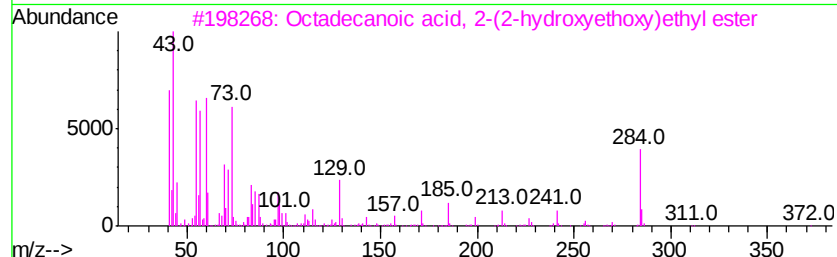
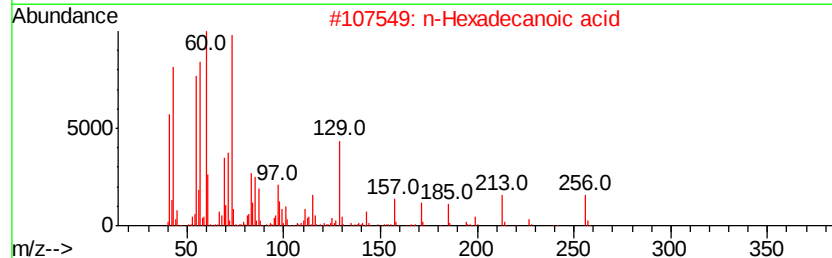
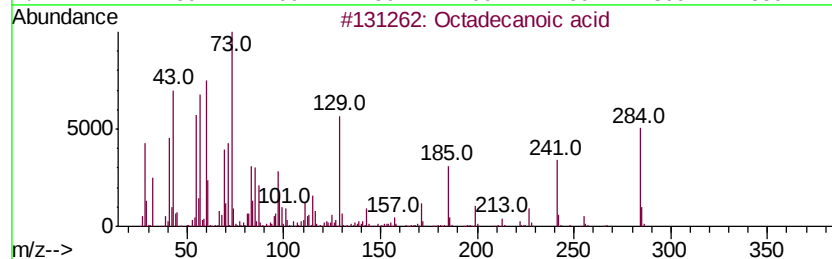
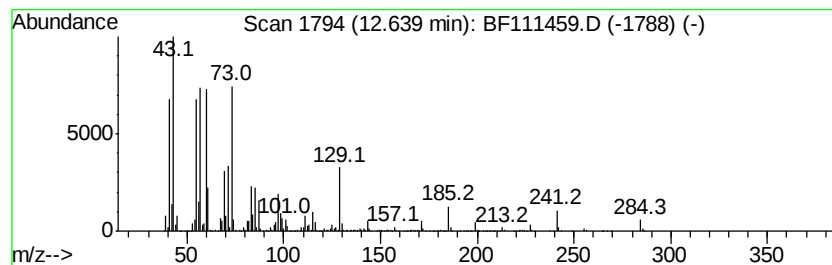
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	78.38 ng	4930060	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
4		Tridecanoic acid	214	C13H26O2	000638-53-9	70
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111459.D
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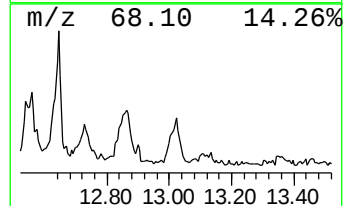
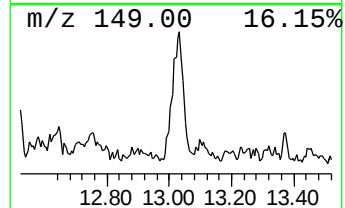
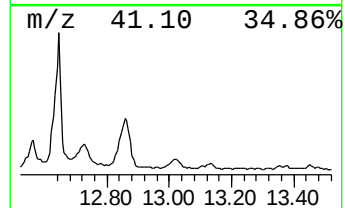
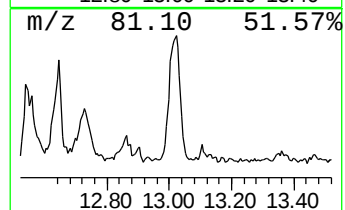
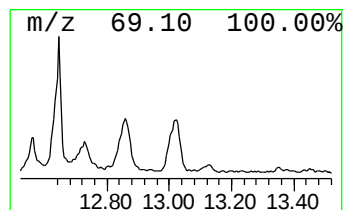
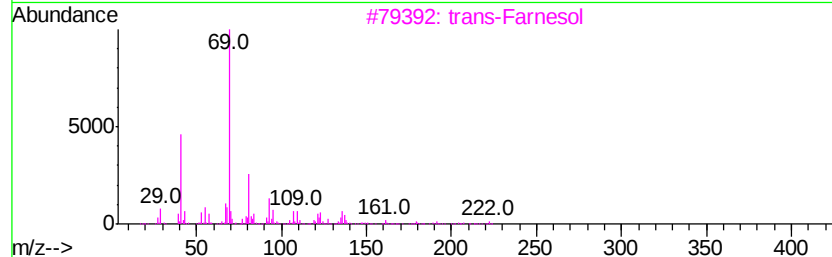
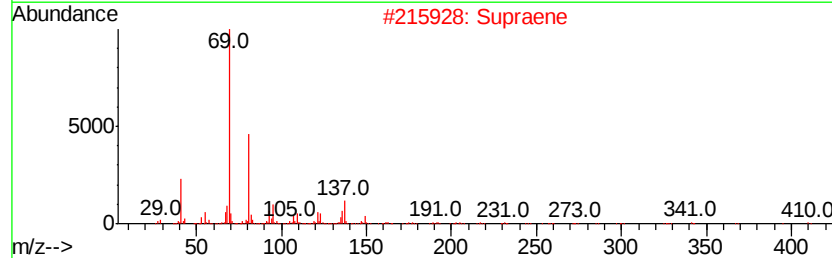
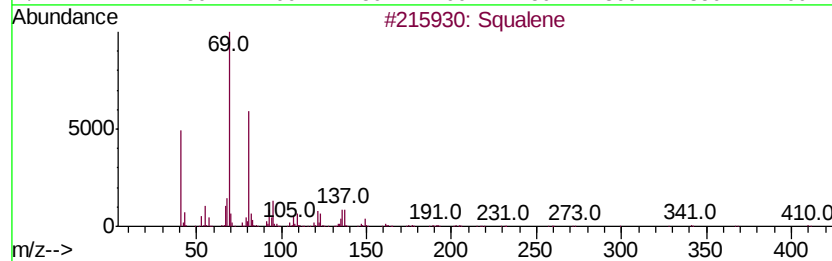
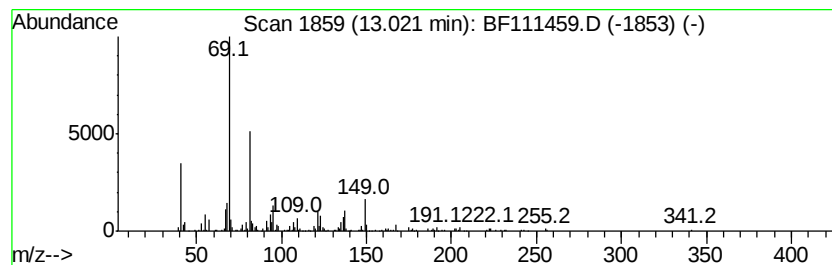
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Squalene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	5.67 ng	356378	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	87
2		Supraene	410	C30H50	007683-64-9	87
3		trans-Farnesol	222	C15H26O	000106-28-5	81
4		Farnesol (E), methyl ether	236	C16H28O	1000352-67-3	64
5		4,8,12-Tetradecatrilal, 5,9,13-...	248	C17H28O	066408-55-7	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
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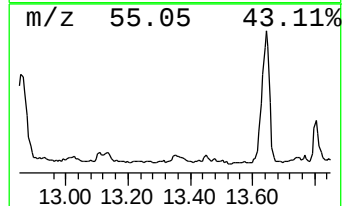
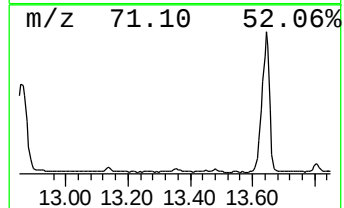
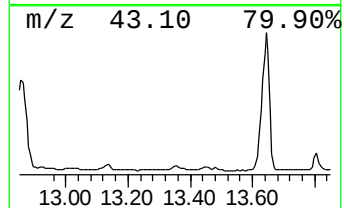
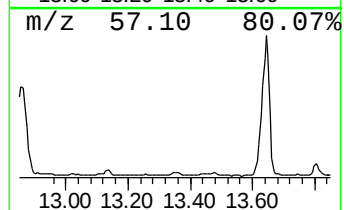
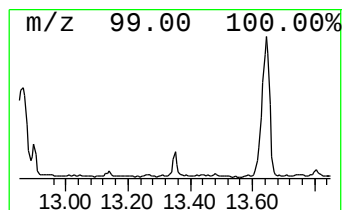
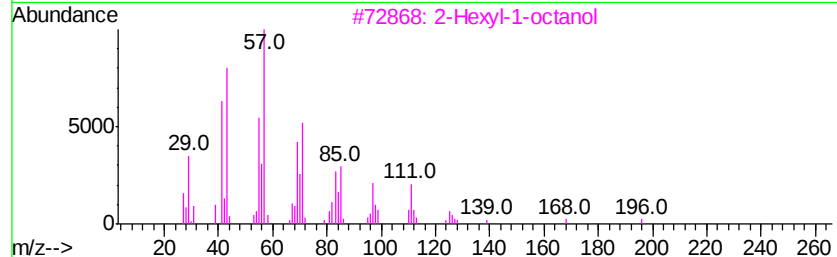
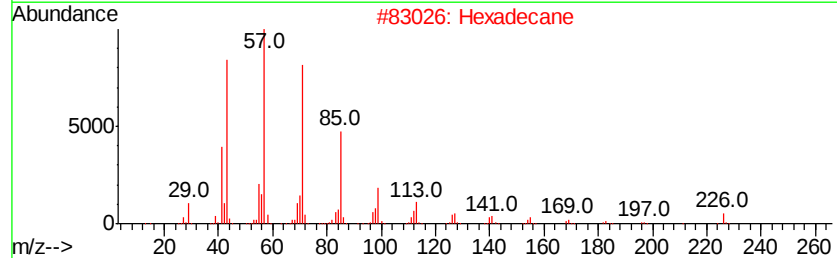
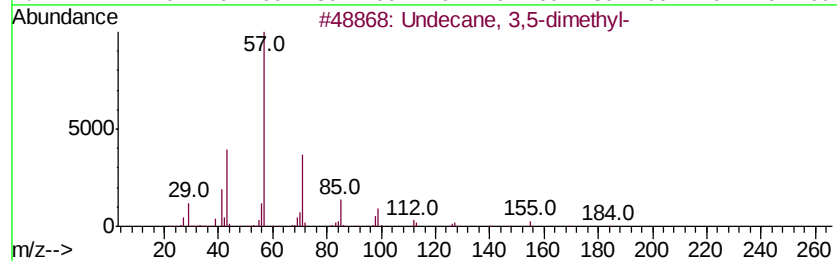
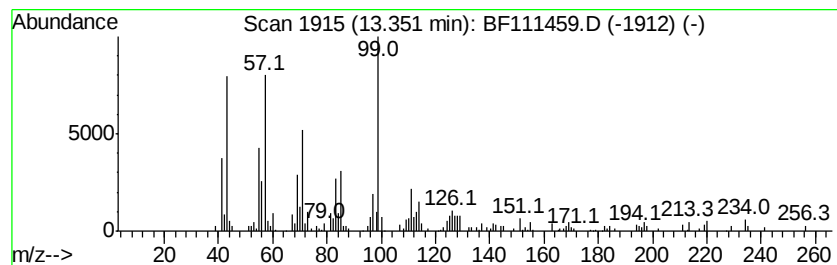
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown13.35 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	3.64 ng	228949	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	41
2			Hexadecane	226	C16H34	000544-76-3	38
3			2-Hexyl-1-octanol	214	C14H30O	019780-79-1	38
4			Decane, 3-methyl-	156	C11H24	013151-34-3	38
5			Heptacosane	380	C27H56	000593-49-7	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
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Sample : J6208-01
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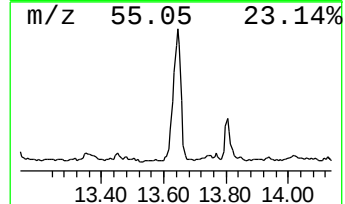
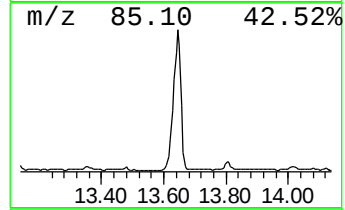
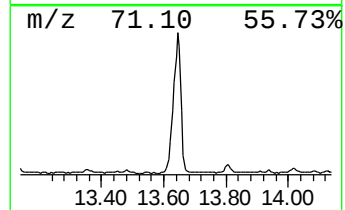
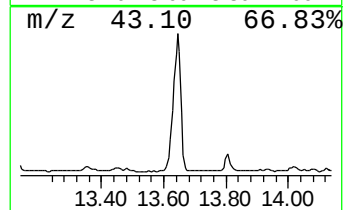
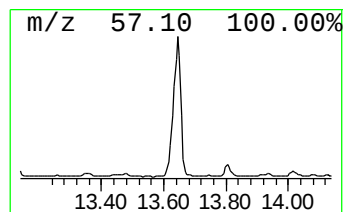
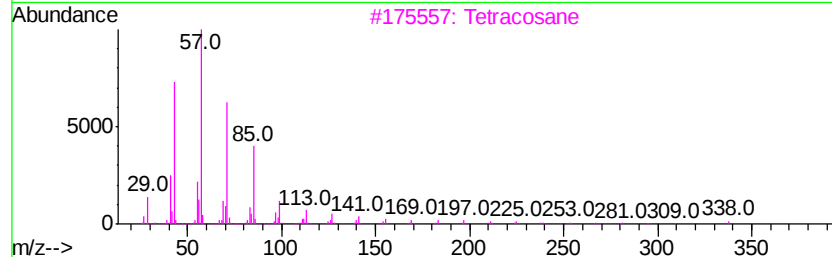
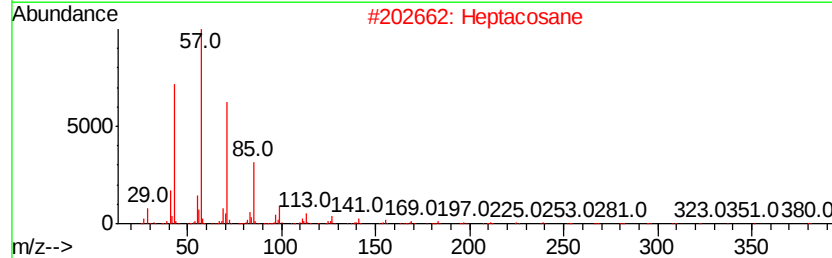
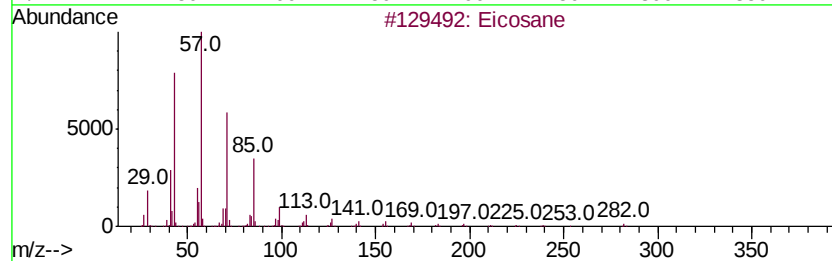
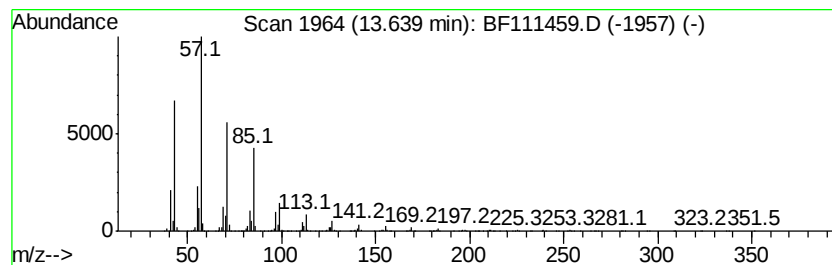
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.64	75.37 ng	4740340	Chrysene-d12		13.95	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Heptacosane		380	C27H56	000593-49-7	91
3	Tetracosane		338	C24H50	000646-31-1	91
4	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	91
5	Hexatriacontane		507	C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
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Instrument :
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ClientSampleId :
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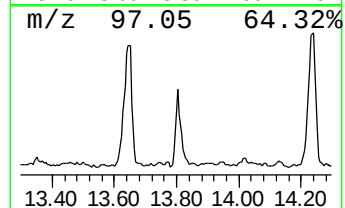
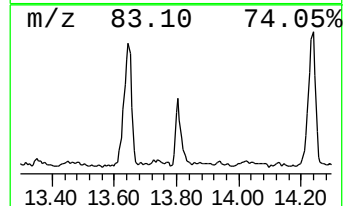
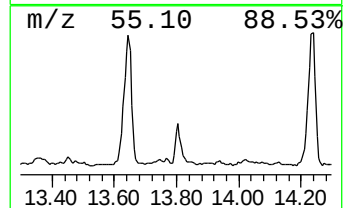
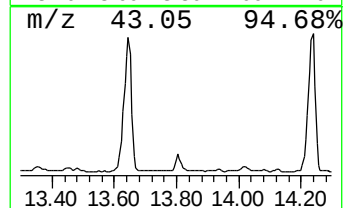
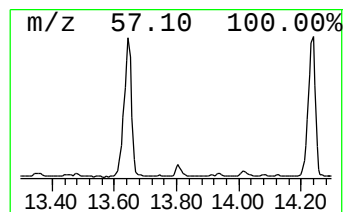
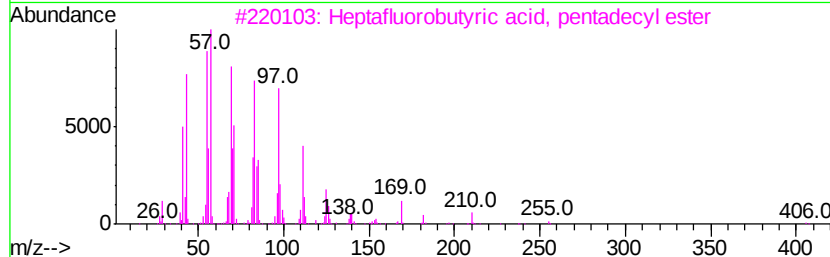
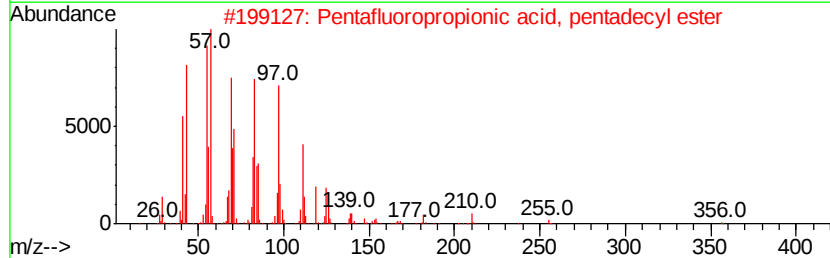
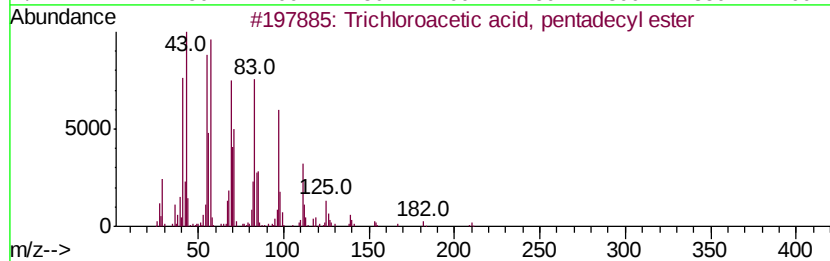
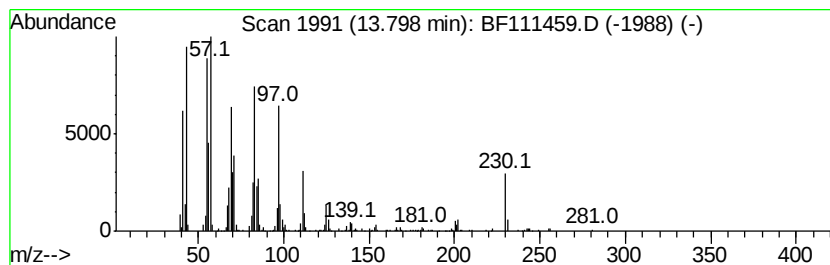
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Trichloroacetic acid, penta... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	11.47 ng	721713	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
2		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
3		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4		1-Heneicosanol	312	C21H44O	015594-90-8	93
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111459.D
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Sample : J6208-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

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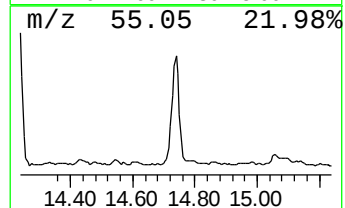
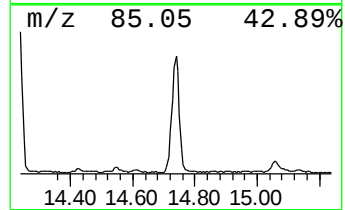
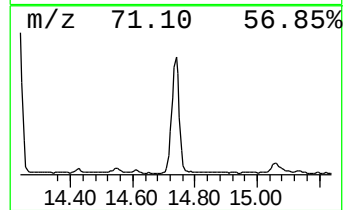
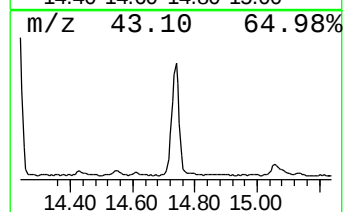
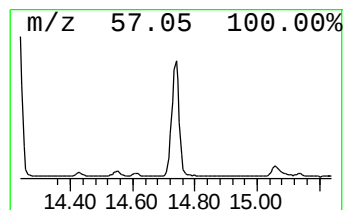
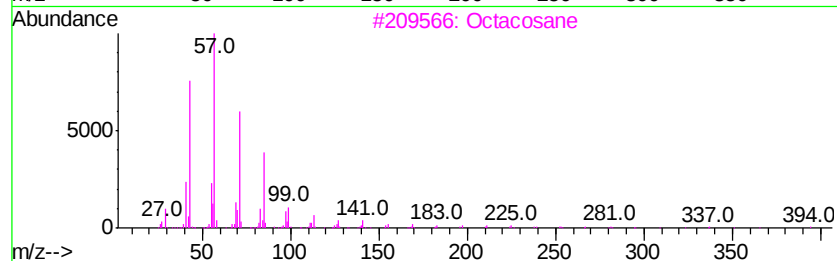
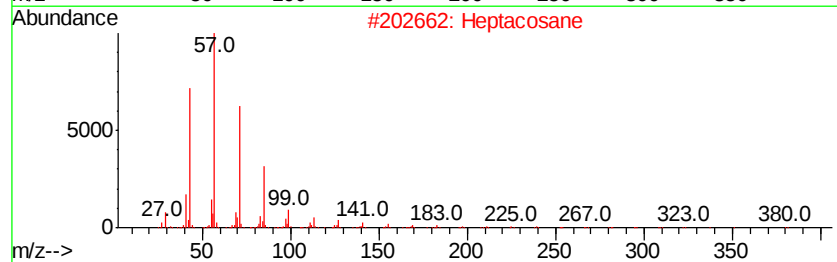
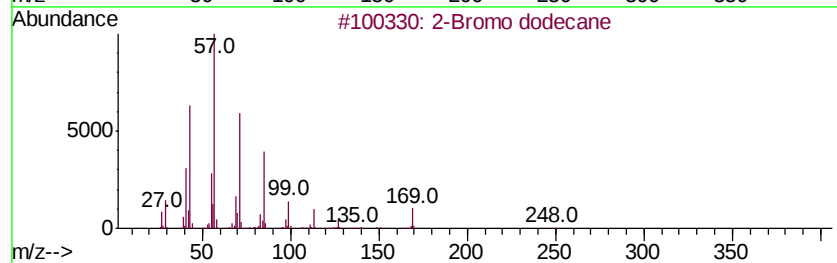
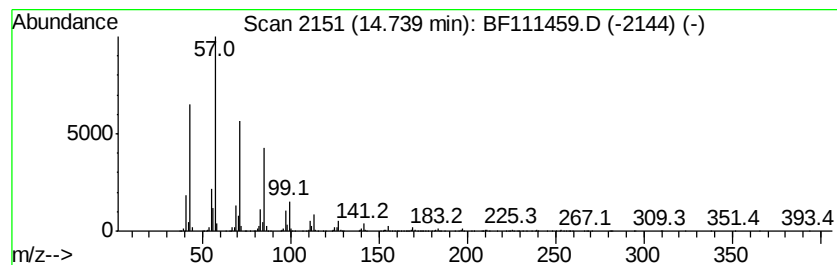
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2-Bromo dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	89.97 ng	3943300	Perylene-d12	15.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
2		Heptacosane	380	C27H56	000593-49-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Pentacosane	352	C25H52	000629-99-2	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
Data File : BF111459.D
Acq On : 15 Dec 2018 10:55
Operator : JU/SJ
Sample : J6208-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-1

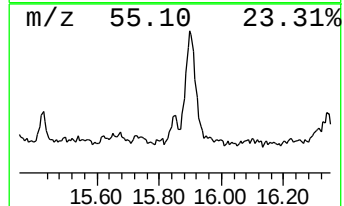
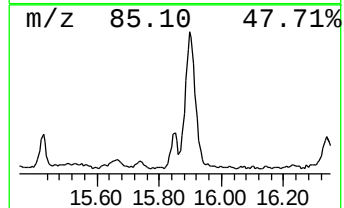
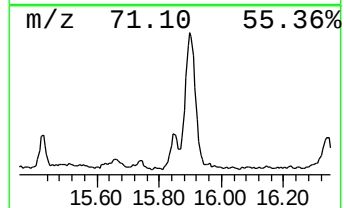
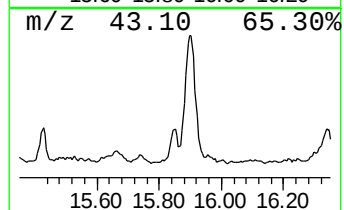
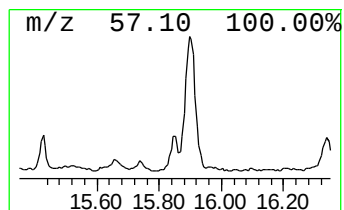
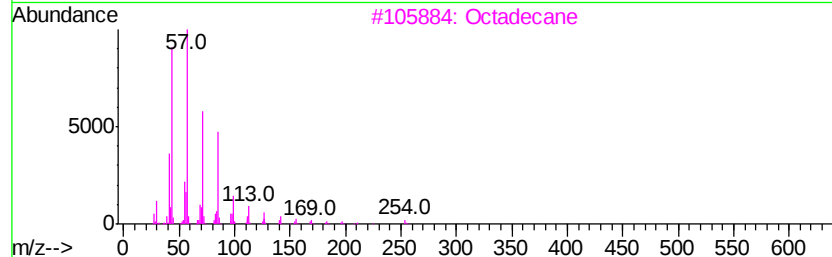
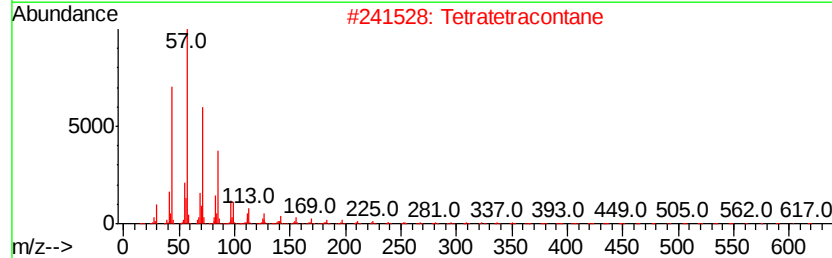
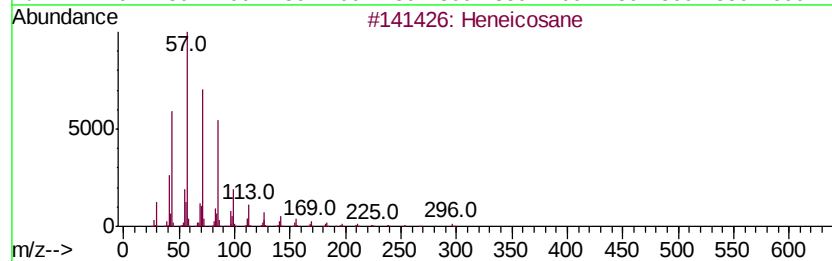
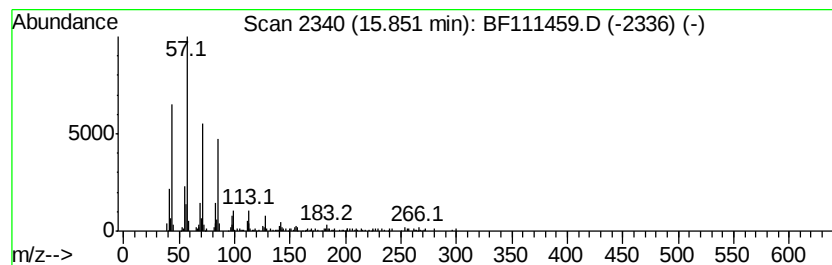
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.85	5.05 ng	221397	Perylene-d12	15.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Tetratetracontane	619	C44H90	007098-22-8	87
3			Octadecane	254	C18H38	000593-45-3	87
4			Octacosane	394	C28H58	000630-02-4	86
5			Hexacosane	366	C26H54	000630-01-3	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121518\
 Data File : BF111459.D
 Acq On : 15 Dec 2018 10:55
 Operator : JU/SJ
 Sample : J6208-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121418.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
unknown-01	3.24	4.1 ng		187986	1	6.79	922588	20.0
2-Pentanone, 4-hy...	5.00	4.8 ng		223120	1	6.79	922588	20.0
unknown6.57	6.57	96.2 ng		4438800	1	6.79	922588	20.0
Tetradecanoic acid	10.99	4.0 ng		265178	4	11.32	1327040	20.0
Octacosane	11.53	20.4 ng		1353430	4	11.32	1327040	20.0
n-Hexadecanoic acid	11.87	131.1 ng		8695450	4	11.32	1327040	20.0
Octadecanoic acid	12.64	78.4 ng		4930060	5	13.95	1257940	20.0
Squalene	13.02	5.7 ng		356378	5	13.95	1257940	20.0
unknown13.35	13.35	3.6 ng		228949	5	13.95	1257940	20.0
Eicosane	13.64	75.4 ng		4740340	5	13.95	1257940	20.0
Trichloroacetic a...	13.80	11.5 ng		721713	5	13.95	1257940	20.0
2-Bromo dodecane	14.74	90.0 ng		3943300	6	15.39	876582	20.0
Heneicosane	15.85	5.0 ng		221397	6	15.39	876582	20.0