

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030818\
 Data File : BF103542.D
 Acq On : 9 Mar 2018 3:51
 Operator : SJ/JU
 Sample : J1817-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 030518-JETT-F-D-B

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:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.122	3	6	15	rVB	84511	136340	3.61%	0.553%
2	3.940	309	315	324	rBV	37880	79863	2.11%	0.324%
3	5.087	507	510	531	rBV	54440	124802	3.30%	0.506%
4	5.492	576	579	605	rBV	502401	1130525	29.90%	4.584%
5	6.510	748	752	770	rBV	290115	784639	20.75%	3.181%
6	6.634	770	773	799	rVB	1849308	2343872	61.99%	9.503%
7	6.863	809	812	828	rBV	715853	856113	22.64%	3.471%
8	7.022	835	839	863	rBV	2561520	3062925	81.00%	12.418%
9	7.434	905	909	933	rBV	2006789	2799426	74.03%	11.350%
10	7.586	933	935	950	rVB	46079	76693	2.03%	0.311%
11	8.151	1027	1031	1056	rBV	650720	1123276	29.71%	4.554%
12	9.222	1209	1213	1222	rBV	3398625	3781351	100.00%	15.331%
13	9.286	1222	1224	1230	rVB	69363	89488	2.37%	0.363%
14	9.622	1274	1281	1287	rBV	128641	183269	4.85%	0.743%
15	9.822	1311	1315	1319	rBV	119825	117248	3.10%	0.475%
16	9.880	1323	1325	1326	rVV	55398	38437	1.02%	0.156%
17	9.904	1326	1329	1342	rVB2	843753	999704	26.44%	4.053%
18	10.316	1396	1399	1402	rVB	144524	112035	2.96%	0.454%
19	10.533	1431	1436	1439	rBV	39144	49307	1.30%	0.200%
20	10.569	1439	1442	1448	rVB	108703	109589	2.90%	0.444%
21	10.704	1461	1465	1477	rBV	843904	1207835	31.94%	4.897%
22	10.792	1477	1480	1489	rVB	123170	163330	4.32%	0.662%
23	11.239	1553	1556	1559	rBV	61359	60281	1.59%	0.244%
24	11.404	1580	1584	1602	rBV	631912	853107	22.56%	3.459%
25	12.986	1848	1853	1863	rBV	2247402	2526347	66.81%	10.243%
26	13.868	1999	2003	2015	rBV2	55258	105463	2.79%	0.428%
27	14.057	2032	2035	2045	rBV	554536	686704	18.16%	2.784%
28	14.798	2158	2161	2167	rVB	37476	40969	1.08%	0.166%
29	14.874	2170	2174	2177	rVB	38782	41167	1.09%	0.167%
30	15.139	2215	2219	2222	rVB	37965	41789	1.11%	0.169%
31	15.521	2280	2284	2287	rBV2	60074	74677	1.97%	0.303%
32	15.568	2287	2292	2302	rVB	278183	469538	12.42%	1.904%
33	15.962	2355	2359	2364	rVB2	58832	82013	2.17%	0.333%
34	16.474	2441	2446	2456	rVB4	60426	121998	3.23%	0.495%

LSC Area Percent Report

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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:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.074	2544	2548	2559	rvB2	47811	101035	2.67%	0.410%
36	17.786	2663	2669	2677	rvB4	34091	89304	2.36%	0.362%

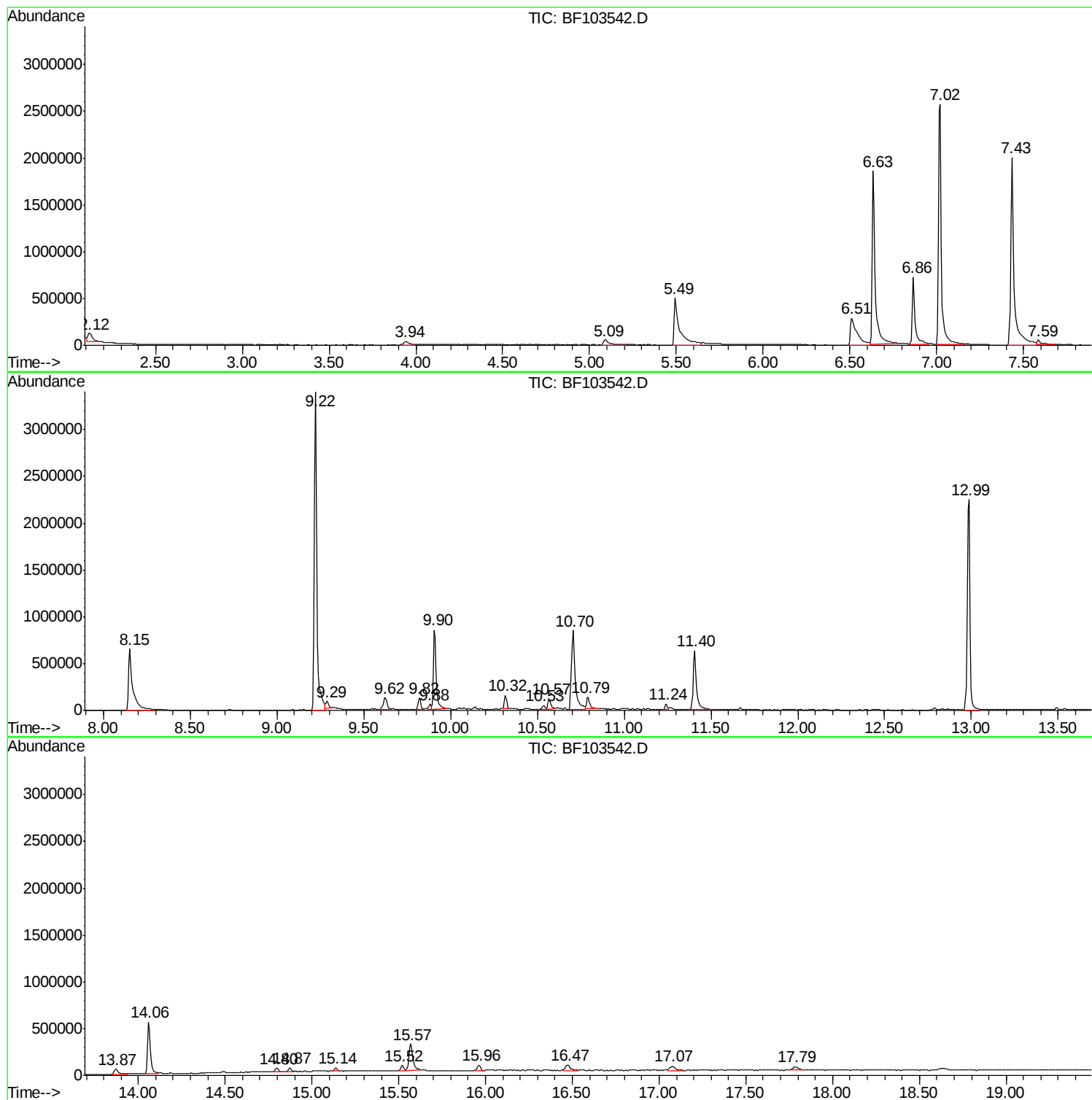
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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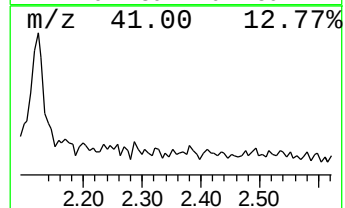
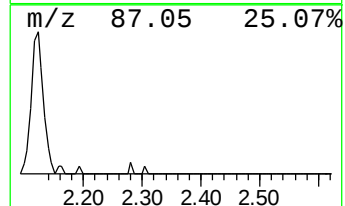
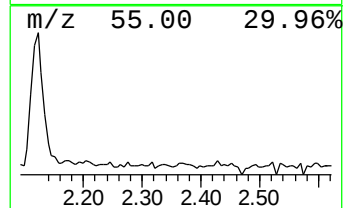
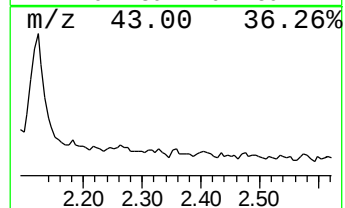
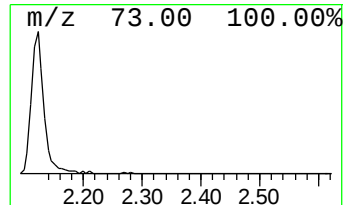
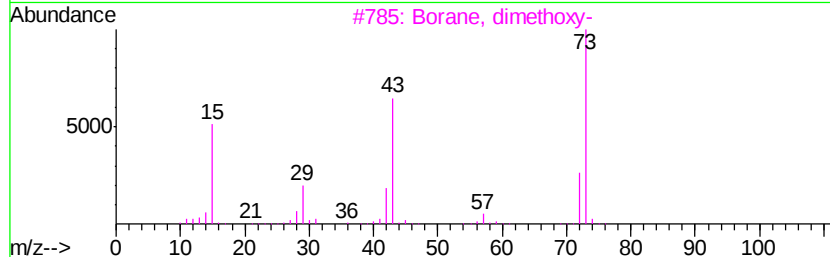
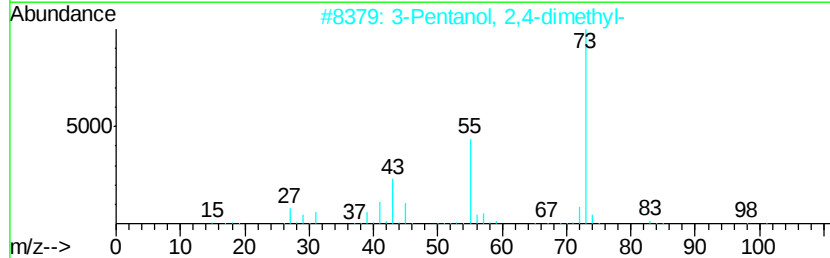
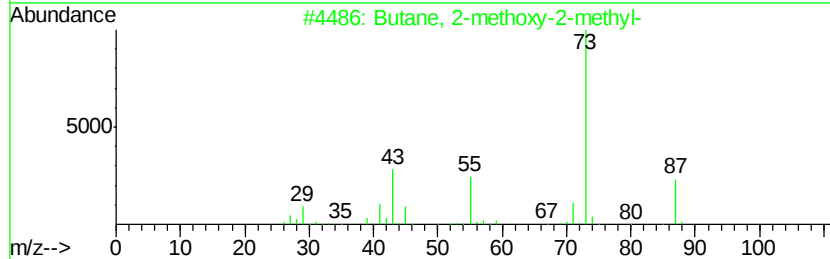
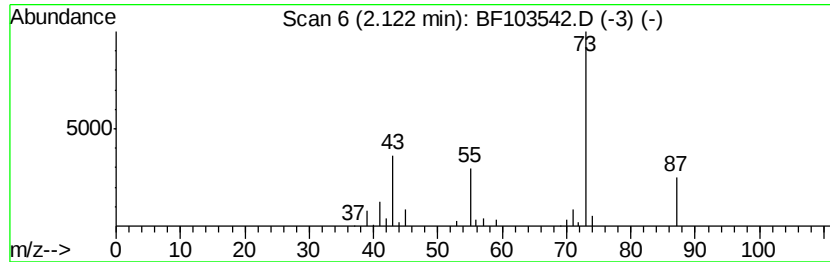
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	3.19 ng	136340	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
2			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	25
3			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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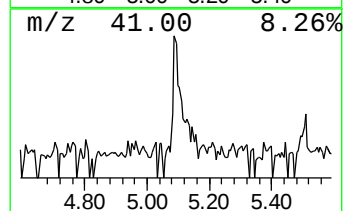
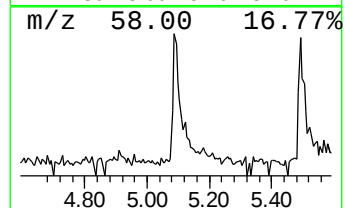
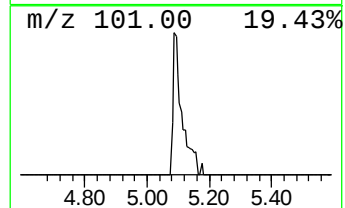
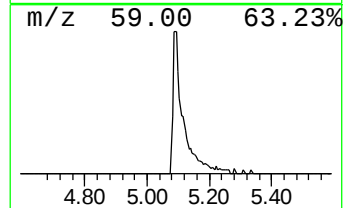
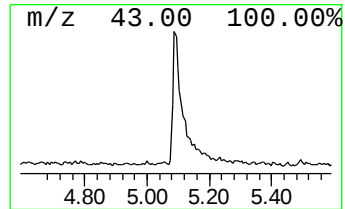
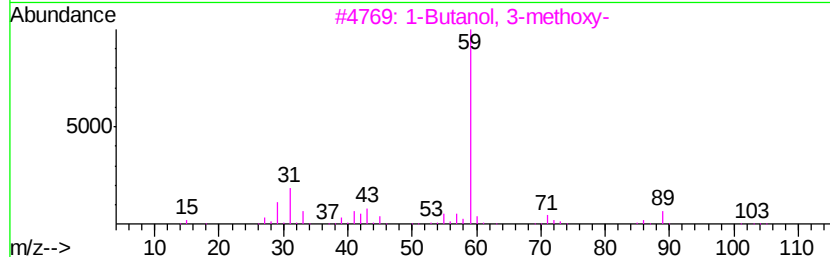
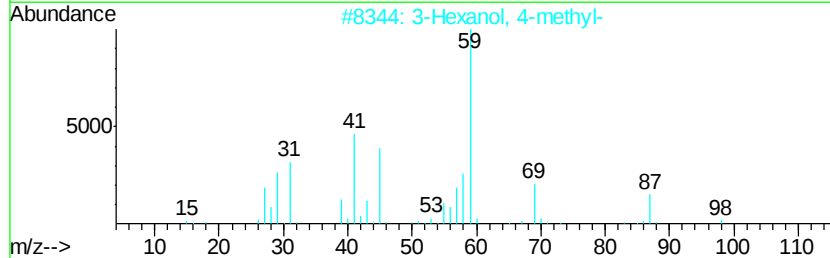
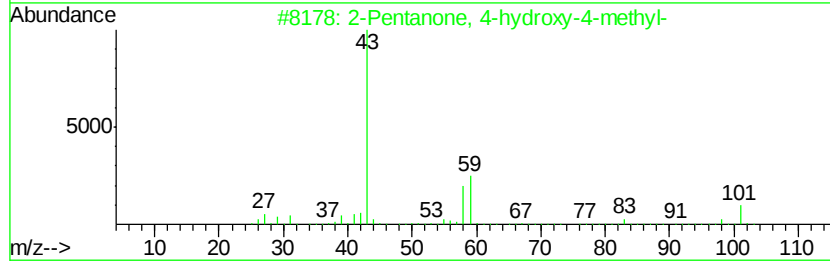
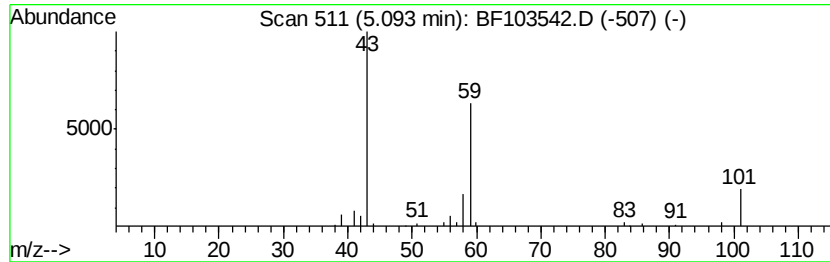
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	2.92 ng	124802	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
3			1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	17
4			4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	9
5			Hexane, 1-ethoxy-	130	C8H18O	005756-43-4	9



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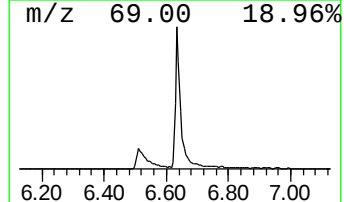
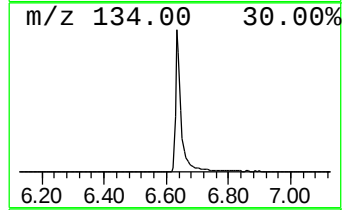
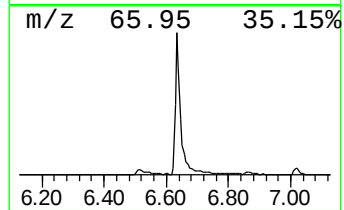
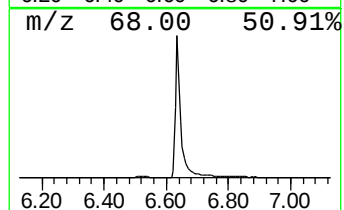
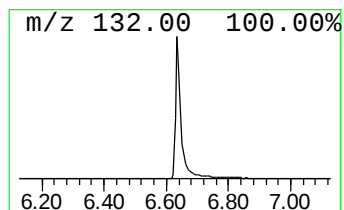
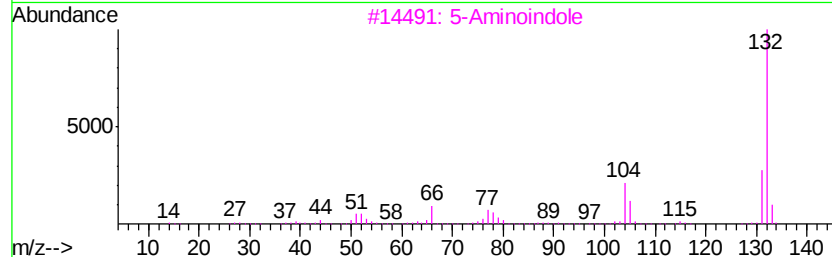
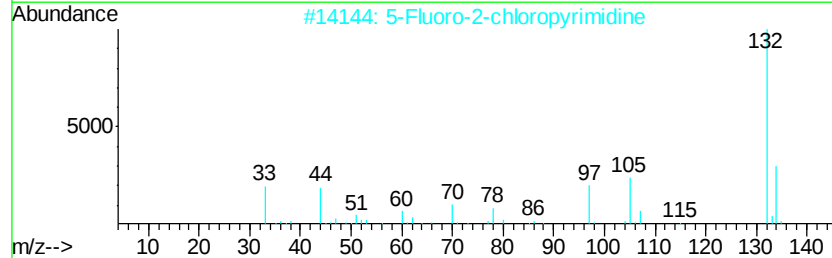
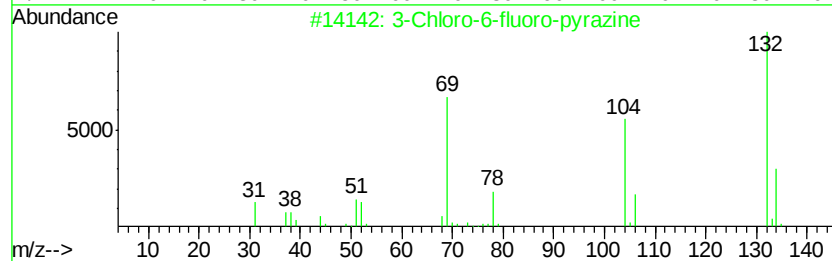
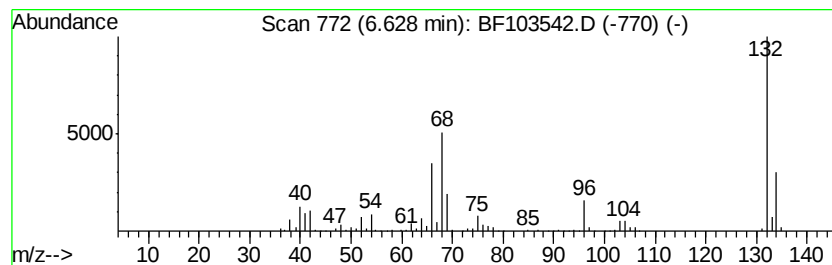
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	54.76 ng	2343870	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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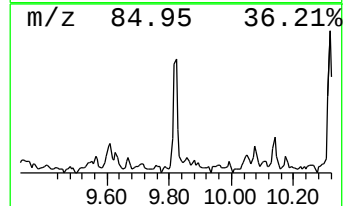
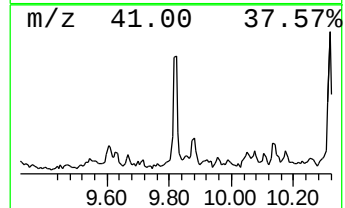
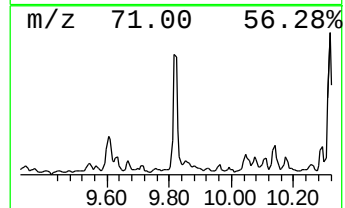
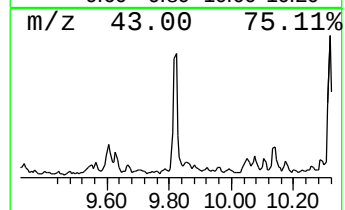
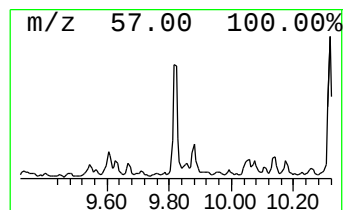
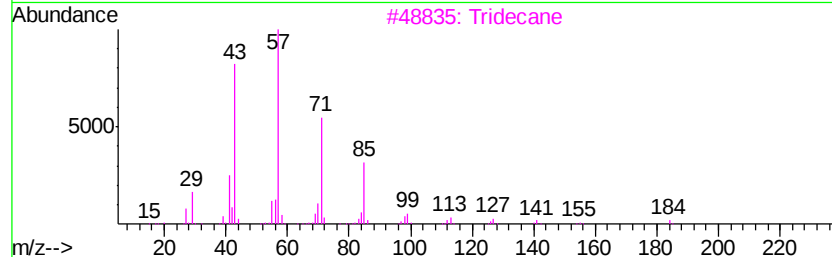
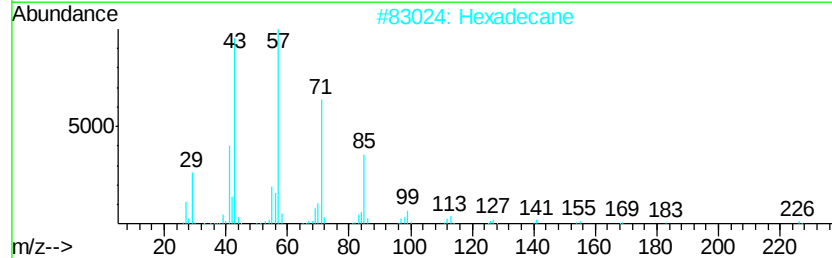
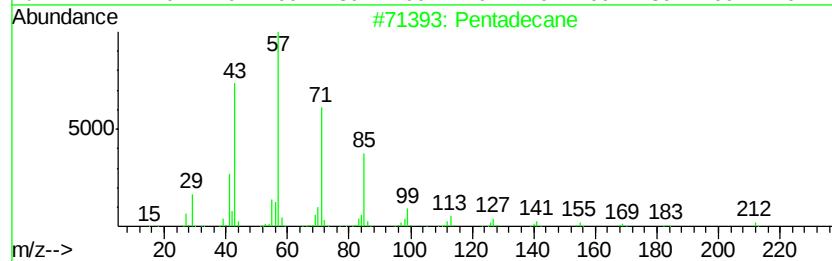
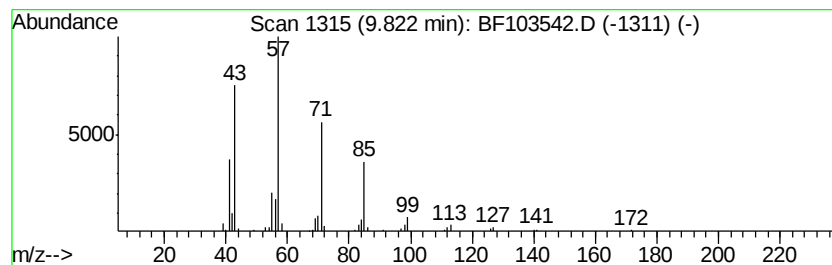
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Pentadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	2.35 ng	117248	Acenaphthene-d10	9.90

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	90
2	Hexadecane		226	C16H34	000544-76-3	90
3	Tridecane		184	C13H28	000629-50-5	86
4	Tetradecane		198	C14H30	000629-59-4	80
5	Undecane, 2-methyl-		170	C12H26	007045-71-8	59



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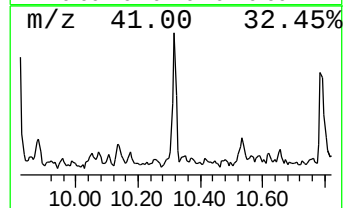
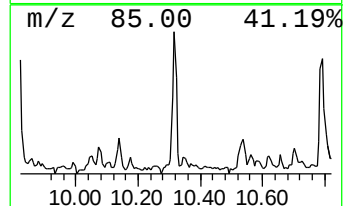
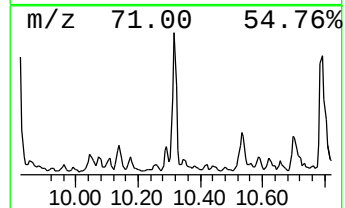
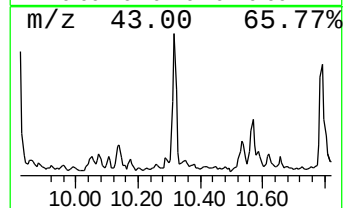
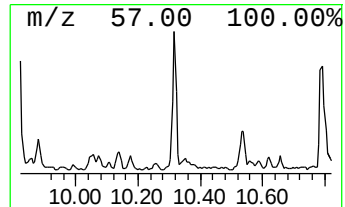
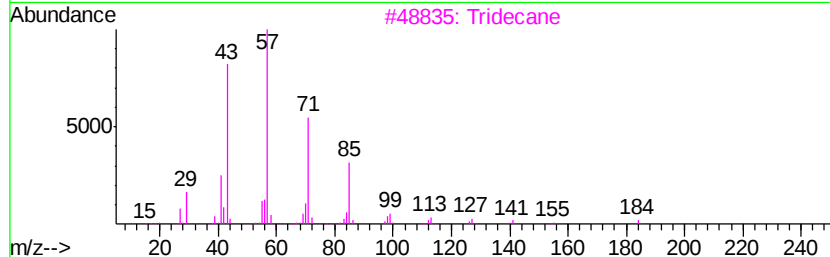
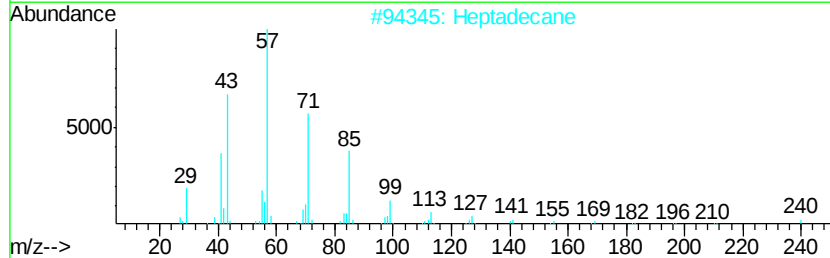
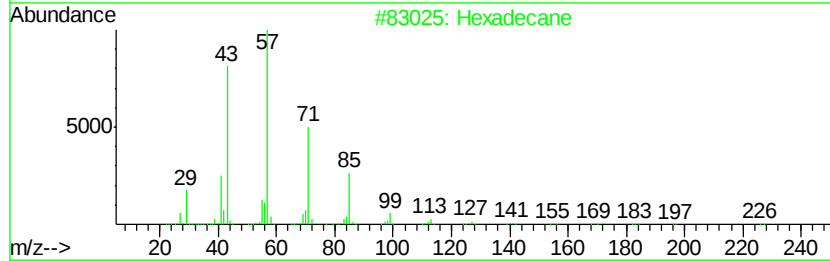
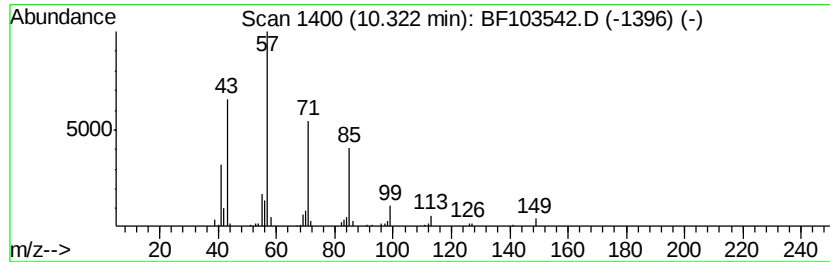
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TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	2.24 ng	112035	Acenaphthene-d10	9.90
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	91
2	Heptadecane	240 C17H36	000629-78-7	86
3	Tridecane	184 C13H28	000629-50-5	86
4	Tetradecane	198 C14H30	000629-59-4	80
5	Sulfurous acid, 2-ethylhexyl hex...	278 C14H30O3S	1000309-20-2	80



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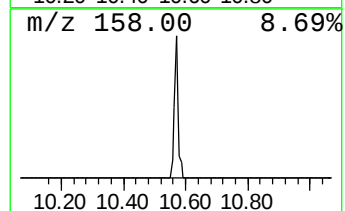
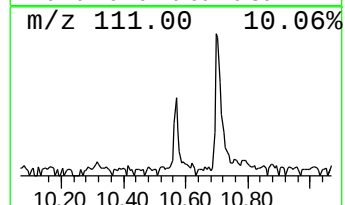
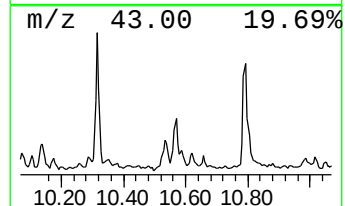
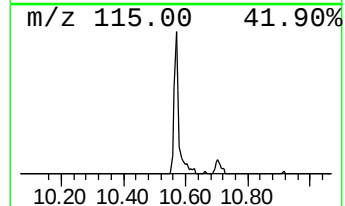
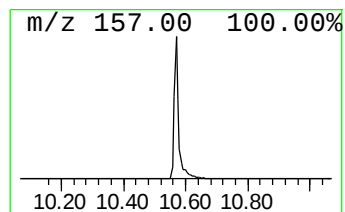
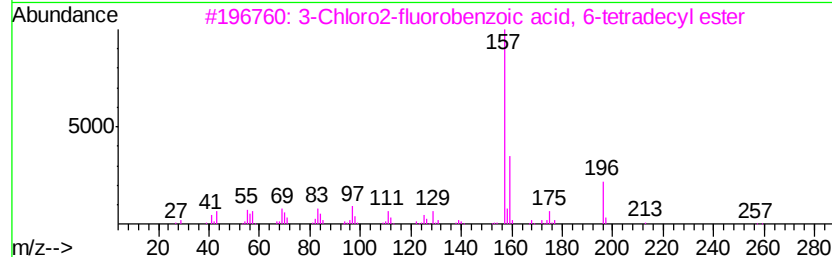
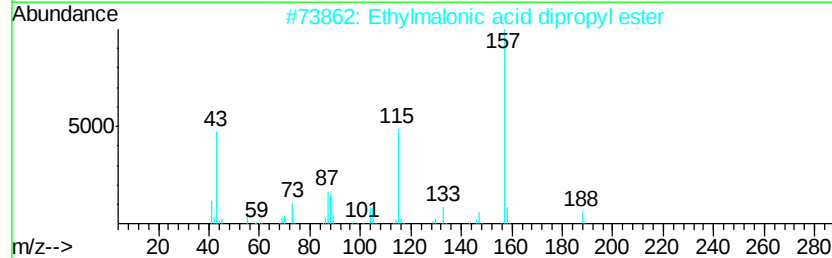
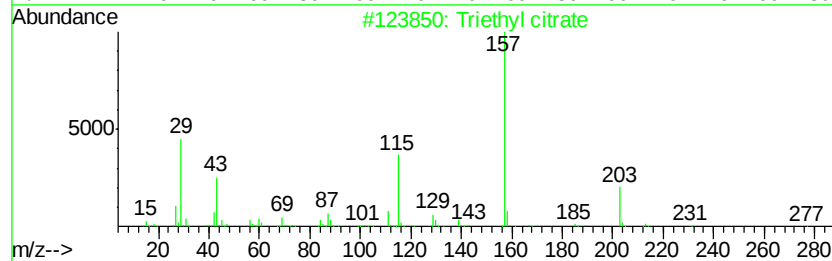
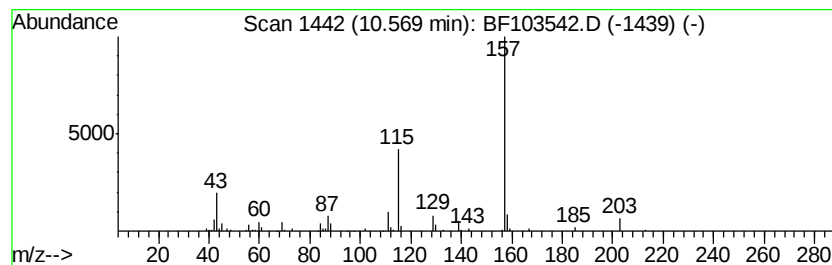
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Triethyl citrate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.57	2.19 ng	109589	Acenaphthene-d10	9.90
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Triethyl citrate	276 C12H20O7	000077-93-0 90
2		Ethylmalonic acid dipropyl ester	216 C11H20O4	001113-91-3 64
3		3-Chloro2-fluorobenzoic acid, 6-...	370 C21H32ClF02	1000338-64-9 40
4		Piperazine-1-carboxylic acid, 4-...	362 C14H19ClN2O5S	1000303-80-2 38
5		N-(5-[3-(2-Oxo-benzooxazol-3-yl)]...	318 C14H14N4O3S	1000300-11-5 38



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030818\
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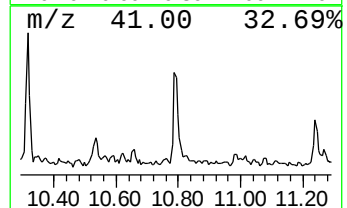
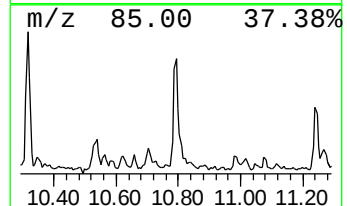
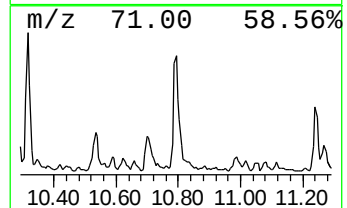
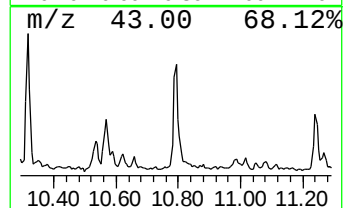
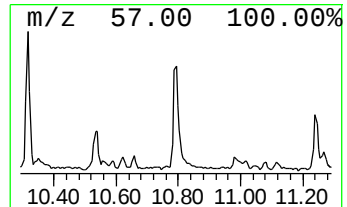
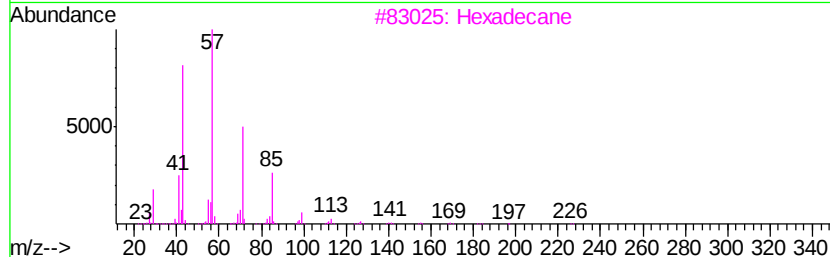
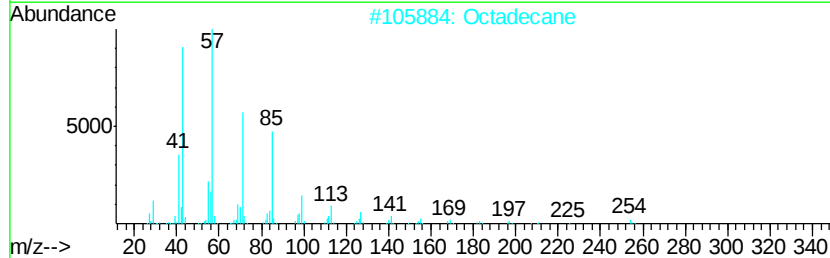
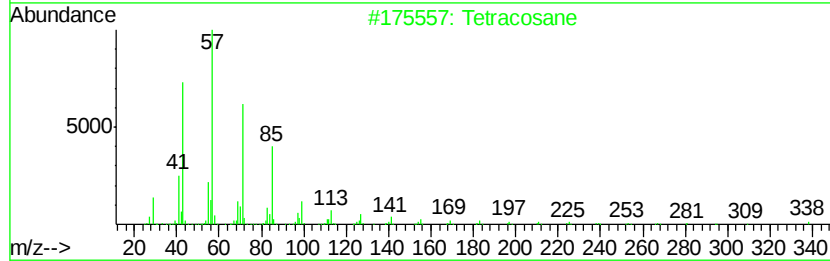
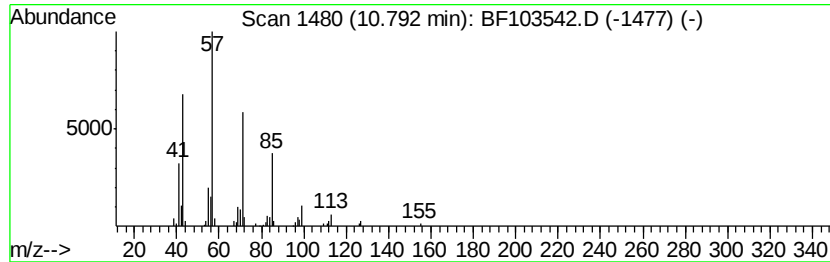
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	3.83 ng	163330	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	90
2		Octadecane	254	C18H38	000593-45-3	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Heneicosane	296	C21H44	000629-94-7	83
5		Heptadecane	240	C17H36	000629-78-7	83



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030818\
Data File : BF103542.D
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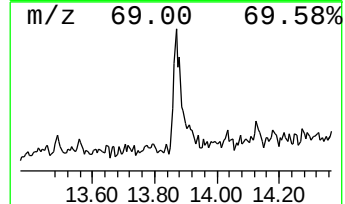
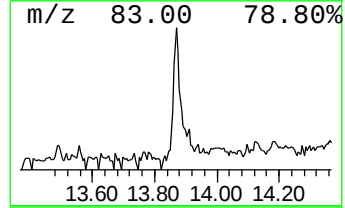
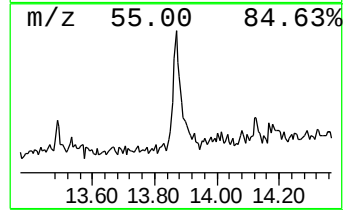
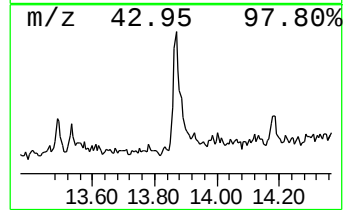
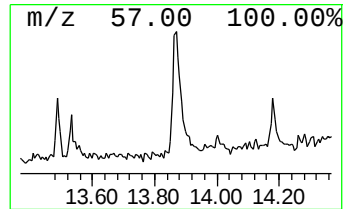
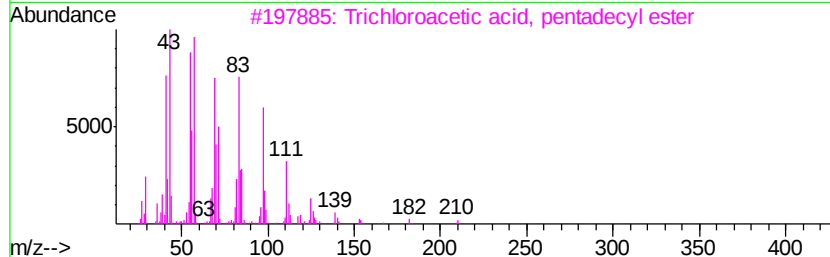
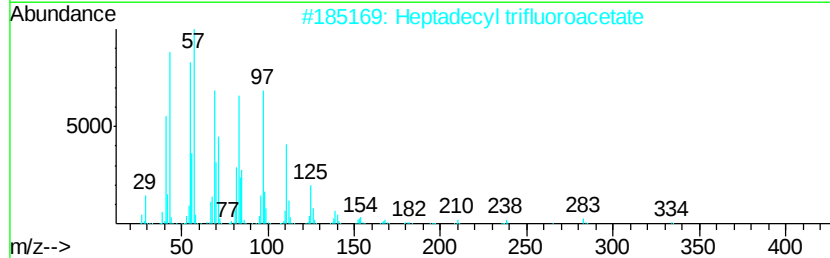
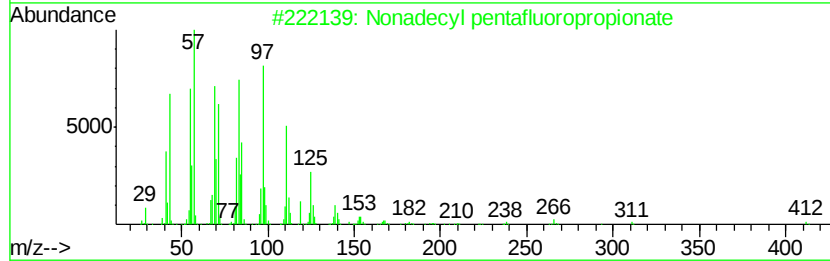
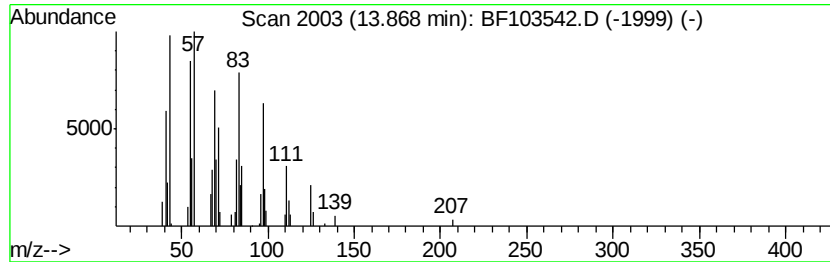
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Nonadecyl pentafluoropropio... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	3.07 ng	105463	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
2		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
3		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	90
5		Octacosanol	410	C28H58O	000557-61-9	90



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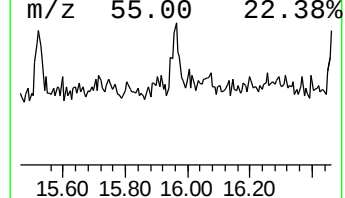
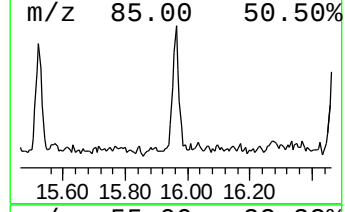
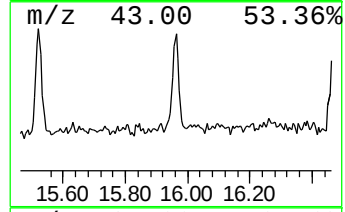
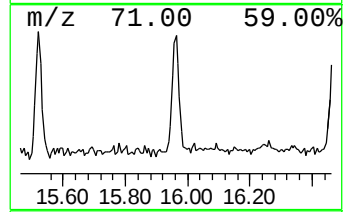
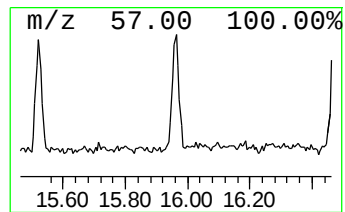
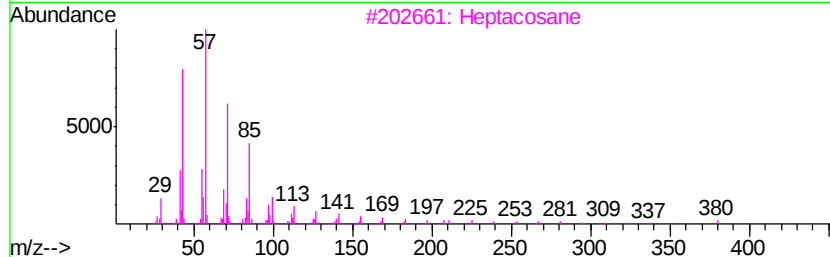
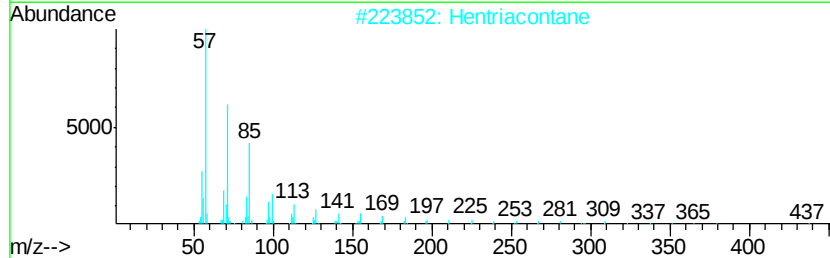
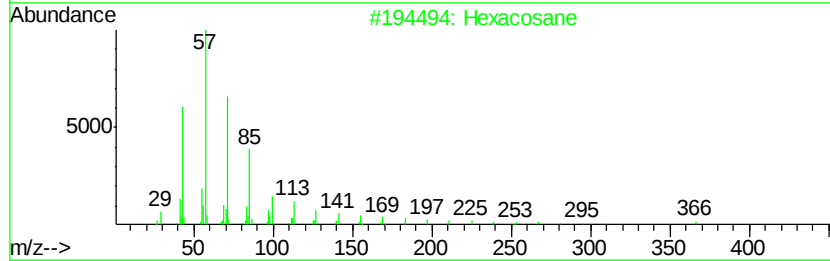
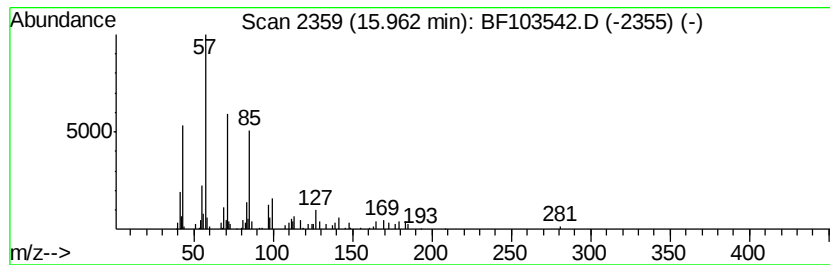
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.96	3.49 ng	82013	Perylene-d12	15.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Hentriacontane	437	C31H64	000630-04-6	90
3	Heptacosane	380	C27H56	000593-49-7	86
4	Heneicosane	296	C21H44	000629-94-7	72
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030818\
Data File : BF103542.D
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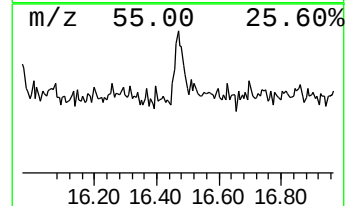
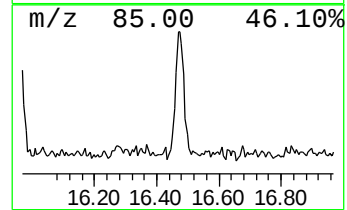
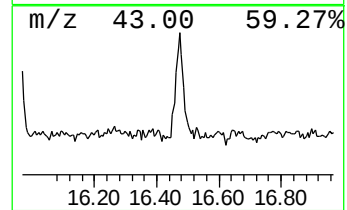
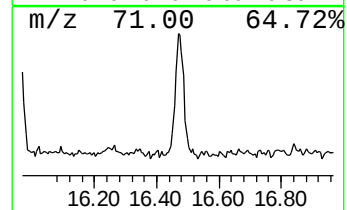
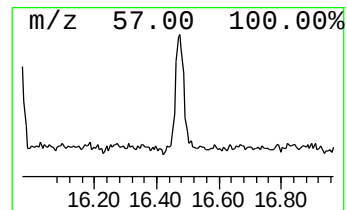
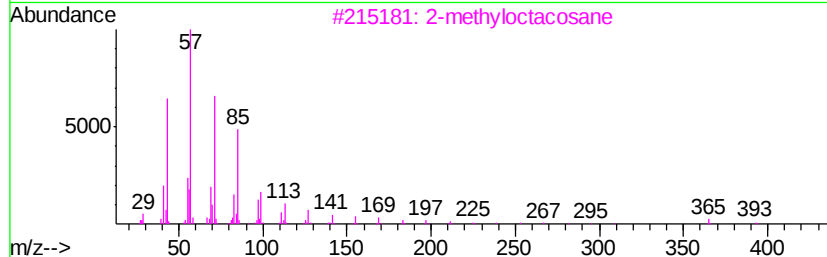
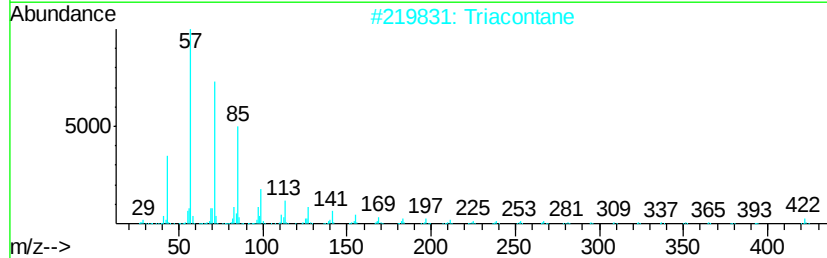
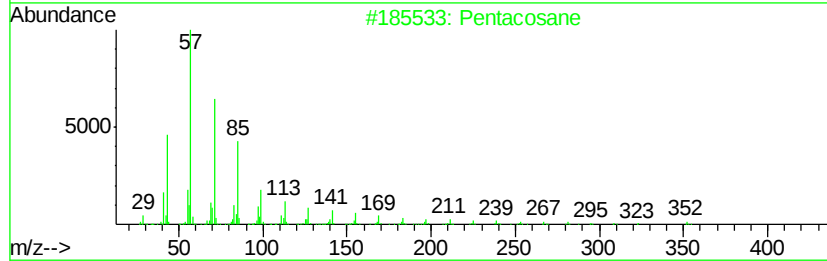
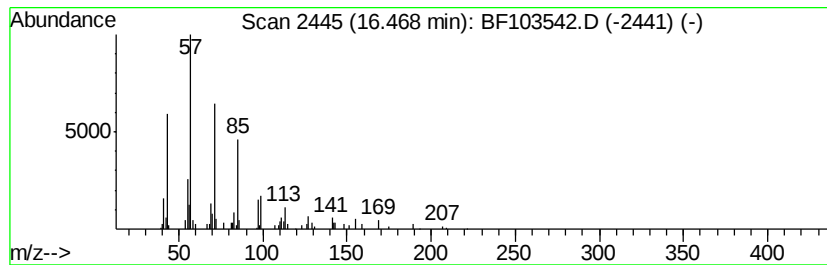
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Pentacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	5.20 ng	121998	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	91
2		Triacontane	422	C30H62	000638-68-6	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	87
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5		2-Bromotetradecane	276	C14H29Br	074036-95-6	86



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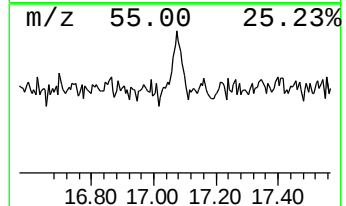
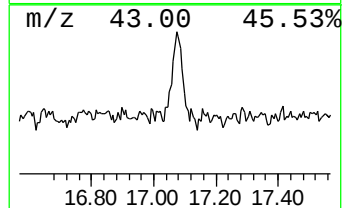
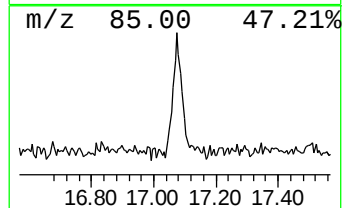
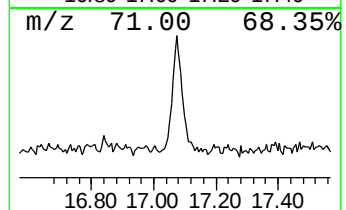
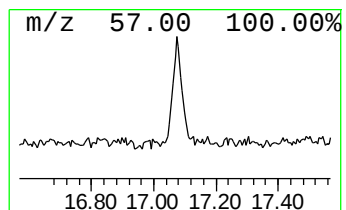
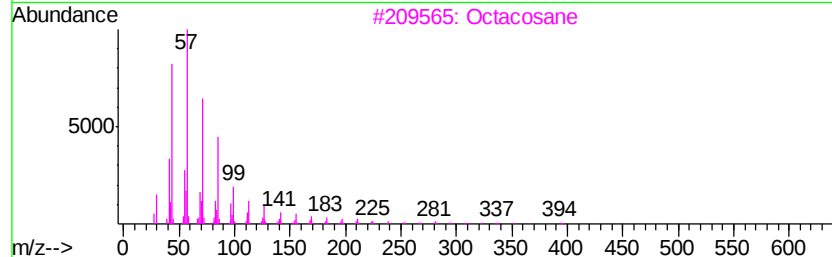
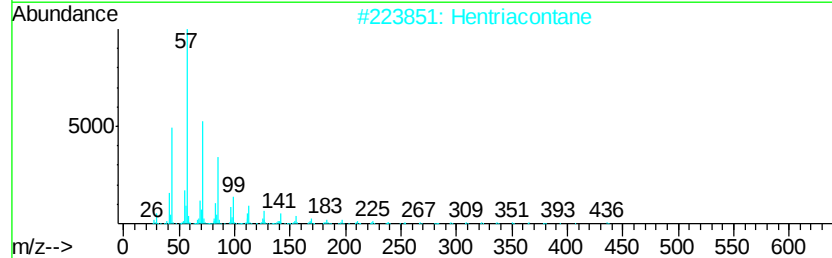
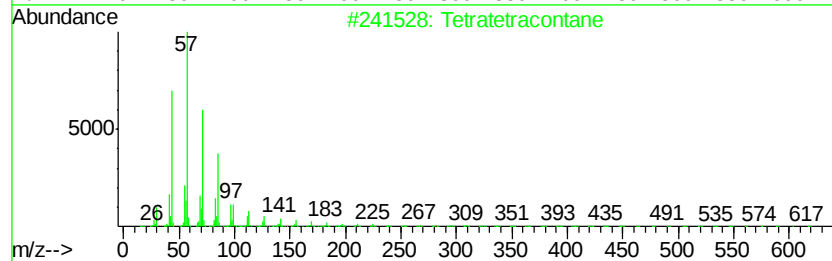
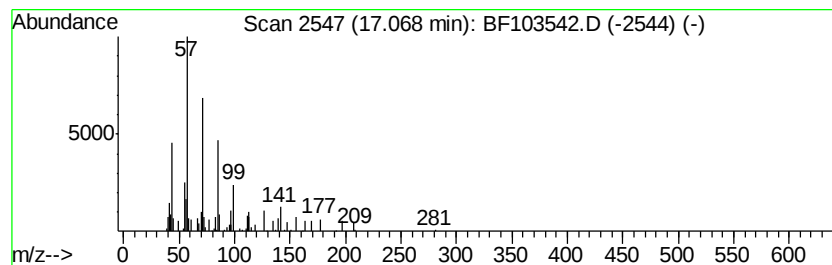
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 12 Tetratetracontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.07	4.30 ng	101035	Perylene-d12	15.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetratetracontane	619	C44H90	007098-22-8	90
2		Hentriacontane	437	C31H64	000630-04-6	87
3		Octacosane	394	C28H58	000630-02-4	87
4		Hexadecane, 2-methyl-	240	C17H36	001560-92-5	86
5		Tetracosane	338	C24H50	000646-31-1	86



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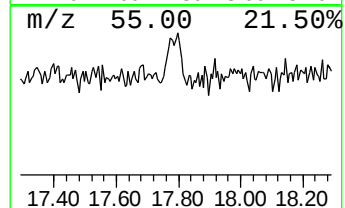
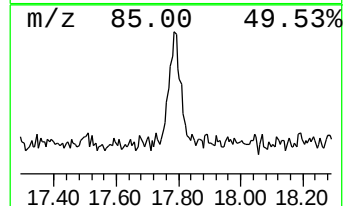
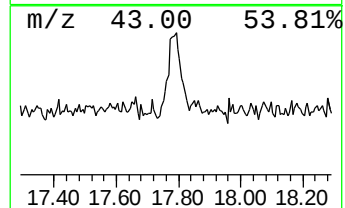
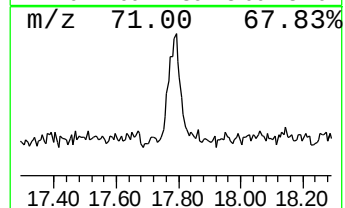
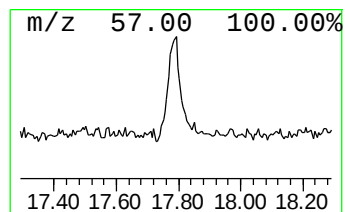
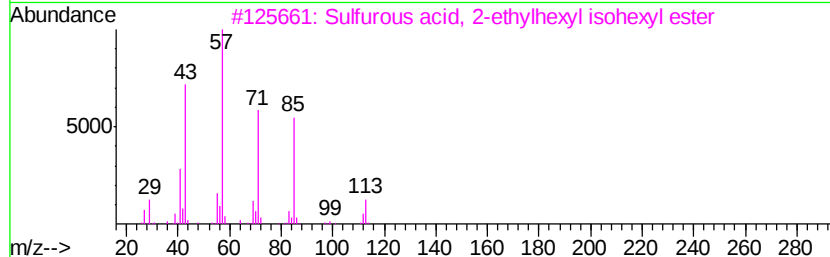
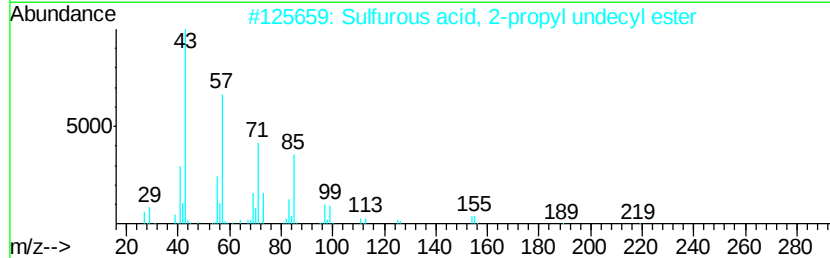
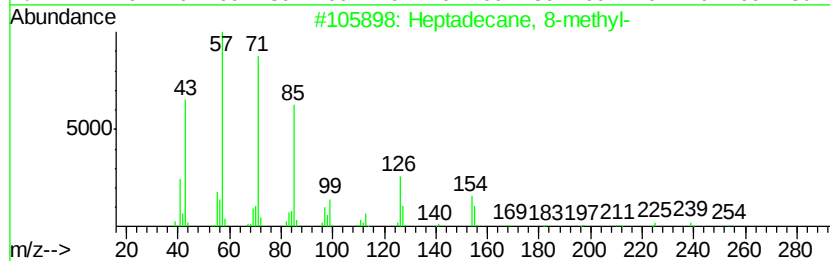
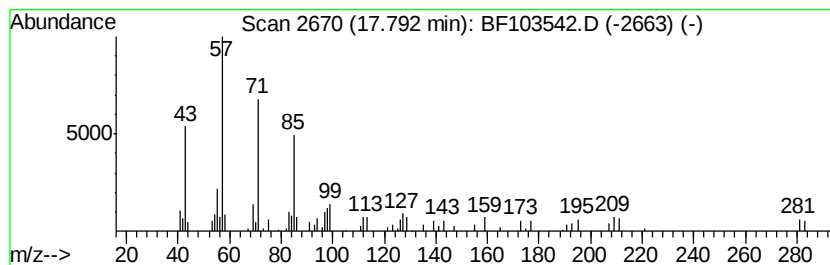
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030518-JETT-F-D-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 8-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.79	3.80 ng	89304	Perylene-d12	15.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane, 8-methyl-	254 C18H38	013287-23-5	80
2	Sulfurous acid, 2-propyl undecyl...	278 C14H30O3S	1000309-12-2	59
3	Sulfurous acid, 2-ethylhexyl iso...	278 C14H30O3S	1000309-19-0	50
4	Sulfurous acid, 2-ethylhexyl hep...	432 C25H52O3S	1000309-20-0	49
5	Sulfurous acid, nonyl 2-propyl e...	250 C12H26O3S	1000309-12-0	47



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030818\
Data File : BF103542.D
Acq On : 9 Mar 2018 3:51
Operator : SJ/JU
Sample : J1817-01
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
030518-JETT-F-D-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.12	3.2	ng	136340	1	6.86	856113	20.0
2-Pentanone, 4-hy...	5.09	2.9	ng	124802	1	6.86	856113	20.0
unknown6.63	6.63	54.8	ng	2343870	1	6.86	856113	20.0
Pentadecane	9.82	2.4	ng	117248	3	9.90	999704	20.0
Hexadecane	10.32	2.2	ng	112035	3	9.90	999704	20.0
Triethyl citrate	10.57	2.2	ng	109589	3	9.90	999704	20.0
Tetracosane	10.79	3.8	ng	163330	4	11.40	853107	20.0
Nonadecyl pentafl...	13.87	3.1	ng	105463	5	14.06	686704	20.0
Hexacosane	15.96	3.5	ng	82013	6	15.57	469538	20.0
Pentacosane	16.47	5.2	ng	121998	6	15.57	469538	20.0
Tetratetracontane	17.07	4.3	ng	101035	6	15.57	469538	20.0
Heptadecane, 8-me...	17.79	3.8	ng	89304	6	15.57	469538	20.0