

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
 Data File : BM017862.D
 Acq On : 01 Dec 2018 18:23
 Operator : JU/SJ
 Sample : J6060-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20180947-COMPOSITE-6

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_M\METHODS\8270-BM111218.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.769	298	303	311	rBV	178359	254436	2.44%	0.287%
2	5.210	372	378	392	rBV	4938205	7012152	67.32%	7.903%
3	6.775	636	644	660	rBV	4532366	7919827	76.04%	8.926%
4	7.122	695	703	715	rBV	5152892	7939795	76.23%	8.948%
5	7.581	773	781	789	rBV	997280	1542925	14.81%	1.739%
6	7.898	827	835	845	rVB	3586020	5646101	54.21%	6.363%
7	8.739	969	978	995	rBV	2825979	4714023	45.26%	5.313%
8	10.351	1245	1252	1273	rBV	1175511	2167967	20.81%	2.443%
9	12.839	1668	1675	1697	rBV	7012163	10358760	99.45%	11.674%
10	13.698	1815	1821	1835	rBV	532563	877018	8.42%	0.988%
11	14.227	1904	1911	1928	rBV2	1812704	2847959	27.34%	3.210%
12	15.727	2159	2166	2179	rBV	5467378	8224429	78.96%	9.269%
13	16.986	2373	2380	2385	rBV	2149266	3088024	29.65%	3.480%
14	17.898	2524	2535	2550	rBV	1098765	1623635	15.59%	1.830%
15	19.056	2725	2732	2740	rBV	142833	283074	2.72%	0.319%
16	19.186	2749	2754	2767	rBV	747650	1081633	10.38%	1.219%
17	19.415	2789	2793	2797	rBV	111840	152905	1.47%	0.172%
18	19.633	2824	2830	2842	rBV	8344719	10415554	100.00%	11.738%
19	20.444	2965	2968	2972	rBV2	114155	140428	1.35%	0.158%
20	20.915	3044	3048	3057	rBV	377178	547364	5.26%	0.617%
21	21.186	3089	3094	3099	rBV	2013919	2640773	25.35%	2.976%
22	21.350	3118	3122	3127	rBV	575833	603518	5.79%	0.680%
23	21.774	3190	3194	3198	rBV	840580	908201	8.72%	1.024%
24	22.227	3266	3271	3276	rBV	894800	1149052	11.03%	1.295%
25	22.715	3349	3354	3360	rBV	964633	1322346	12.70%	1.490%
26	23.262	3442	3447	3456	rVB	708809	1217695	11.69%	1.372%
27	23.368	3459	3465	3481	rVB2	1330937	2601479	24.98%	2.932%
28	23.880	3547	3552	3560	rVB	532437	955456	9.17%	1.077%
29	24.591	3668	3673	3682	rVB	245303	495056	4.75%	0.558%

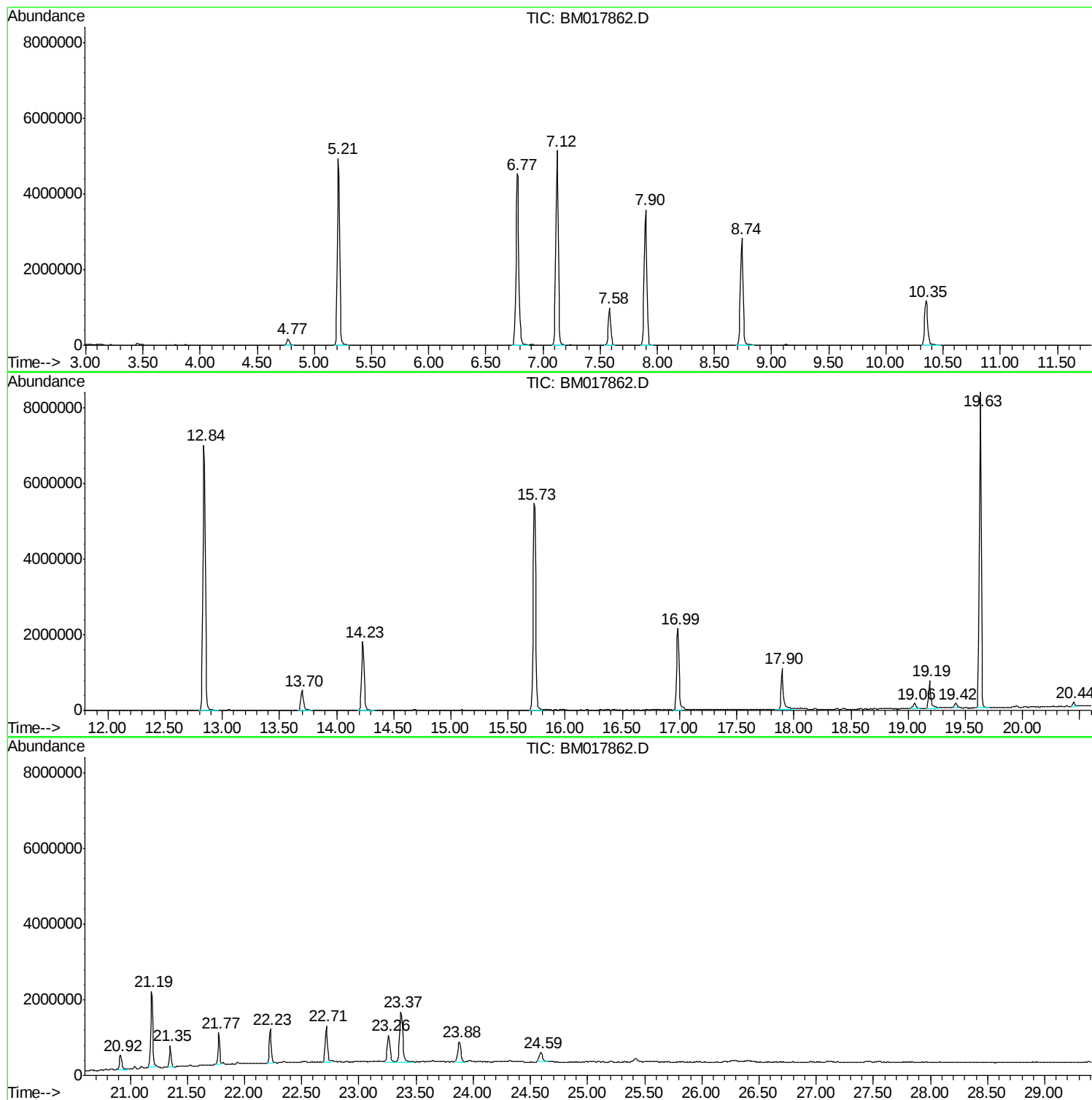
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ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.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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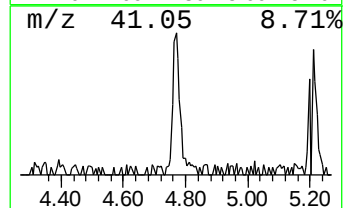
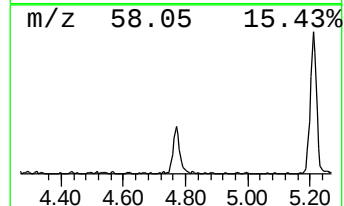
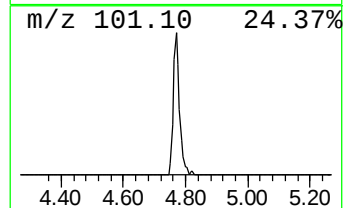
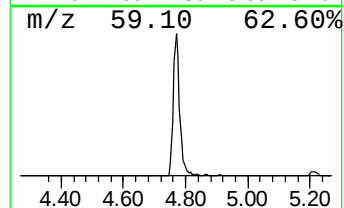
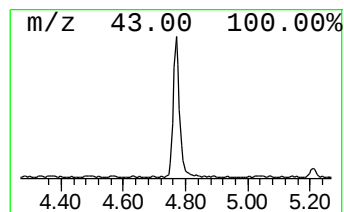
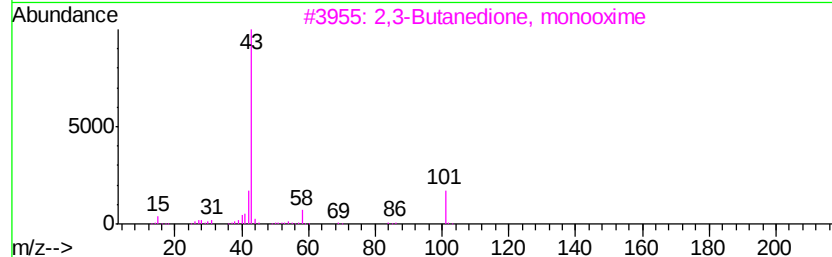
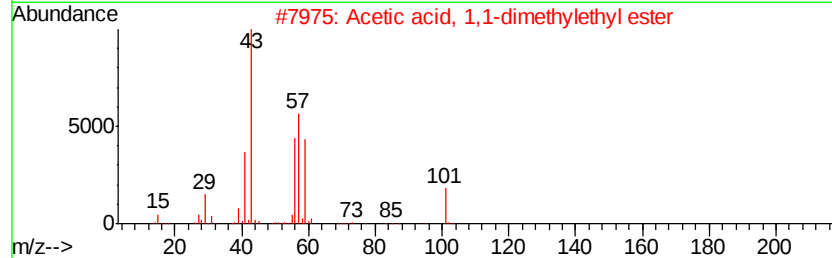
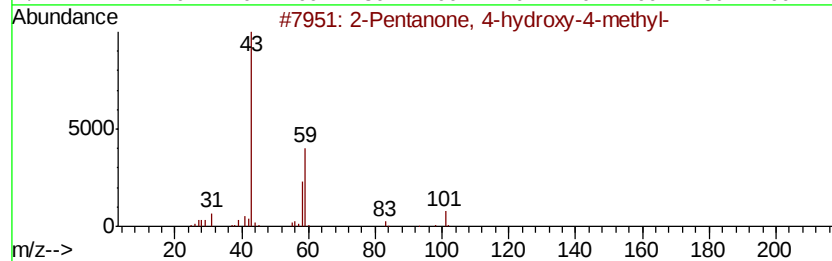
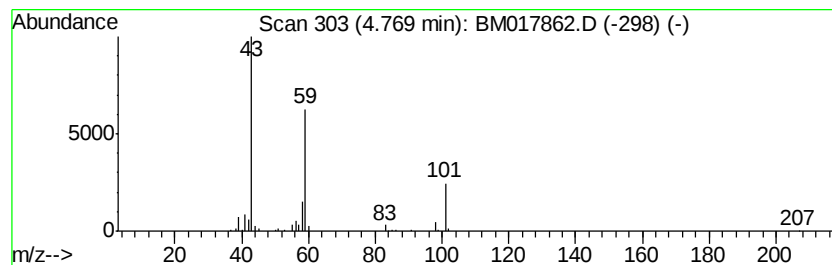
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.77	3.30 ng	254436	1,4-Dichlorobenzene-d4	7.58

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	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
4			1,2-Butanediol, (.+/-.)-	90	C4H10O2	026171-83-5	23
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17



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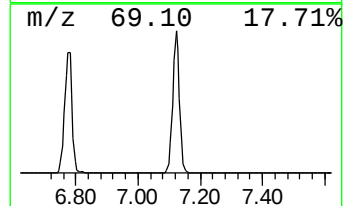
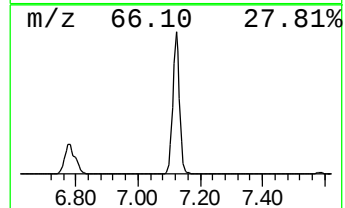
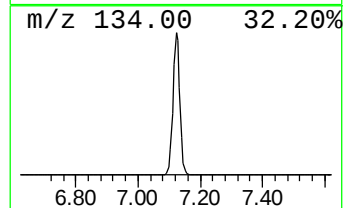
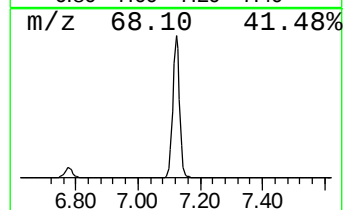
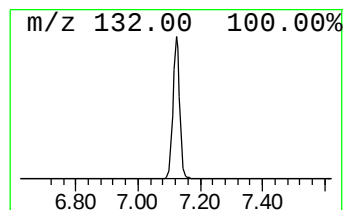
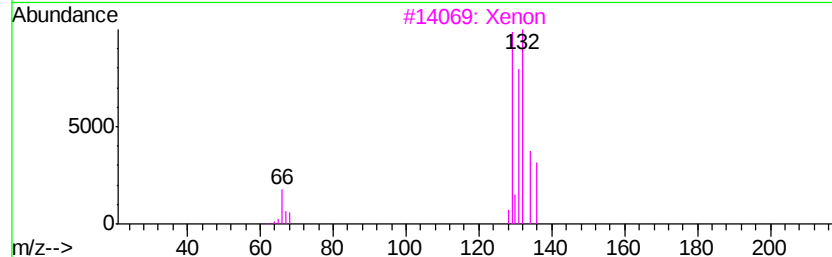
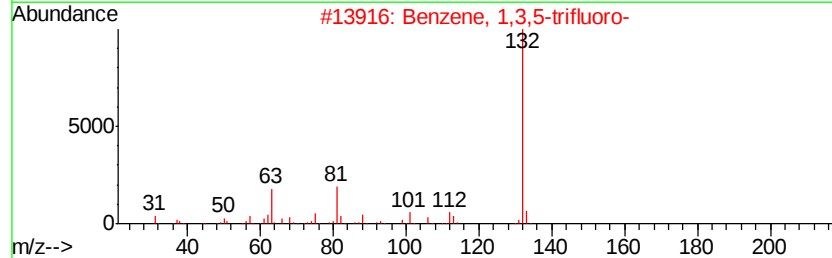
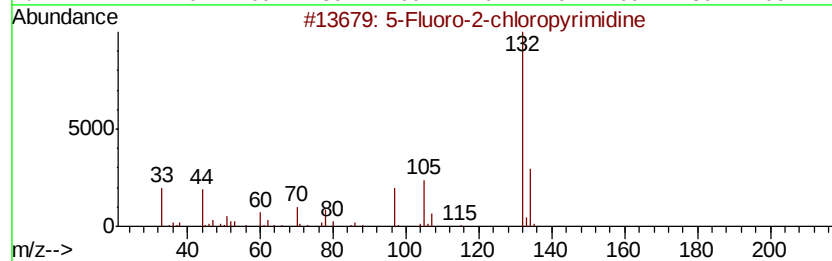
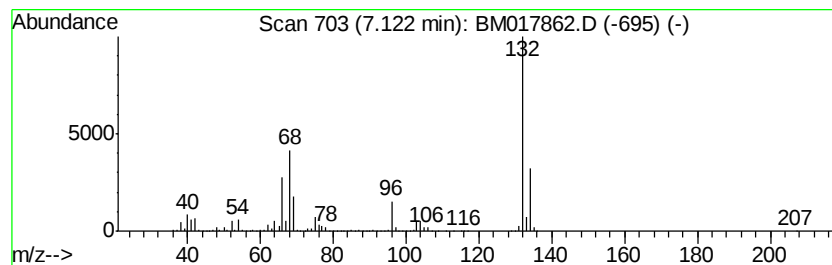
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Peak Number 2 unknown7.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	102.92 ng	7939800	1,4-Dichlorobenzene-d4	7.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
3		Xenon	132	Xe	007440-63-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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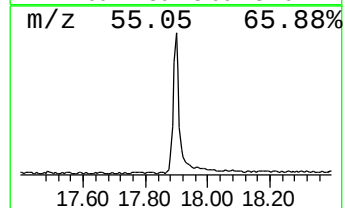
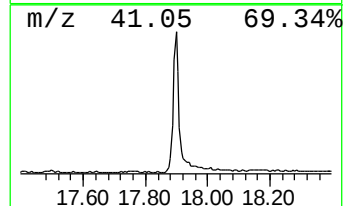
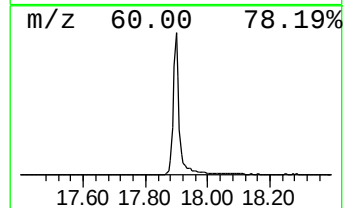
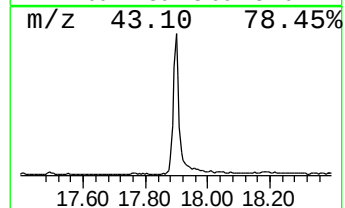
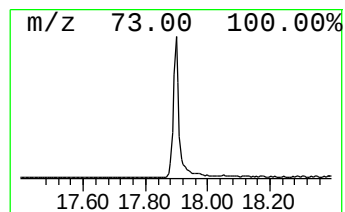
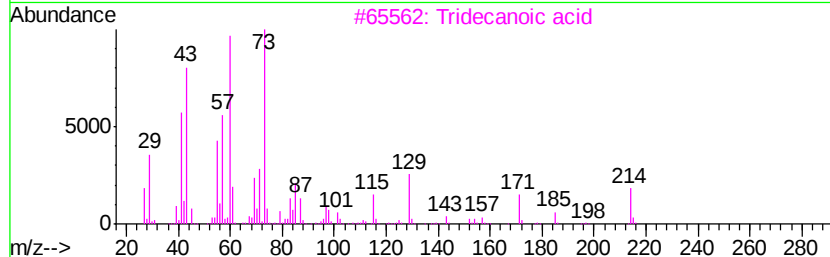
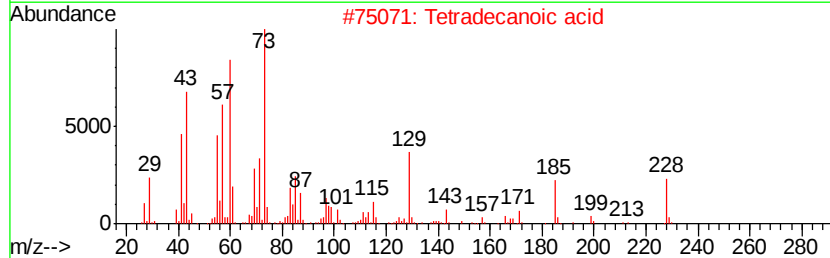
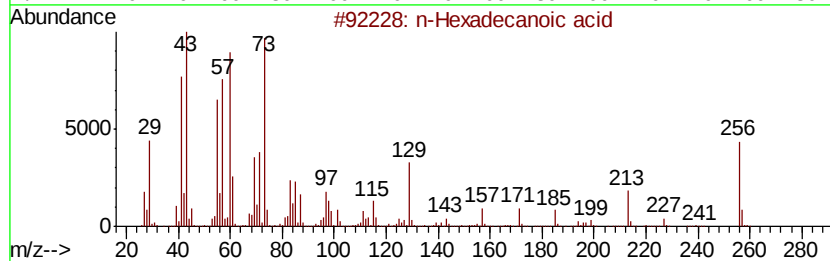
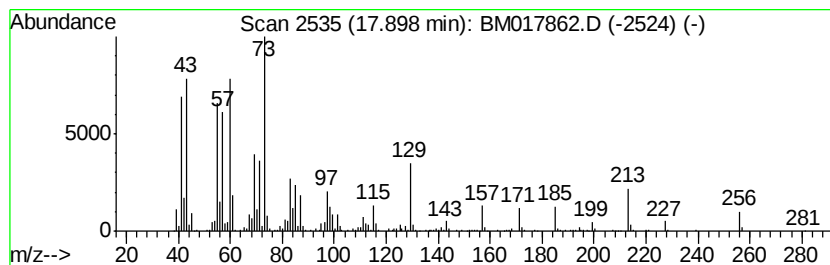
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.90	10.52 ng	1623640	Phenanthrene-d10	16.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	89
3	Tridecanoic acid	214	C13H26O2	000638-53-9	86
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
5	11-Bromoundecanoic acid	264	C11H21BrO2	002834-05-1	47



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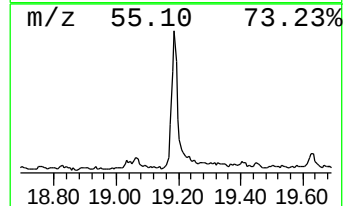
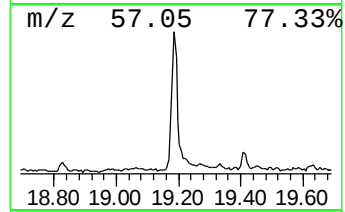
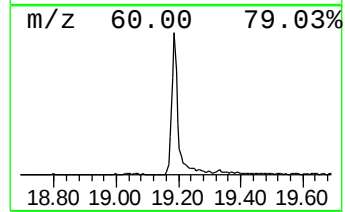
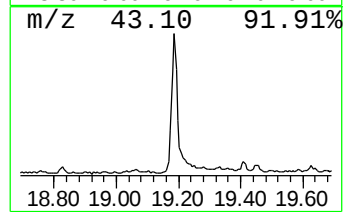
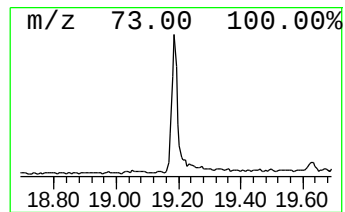
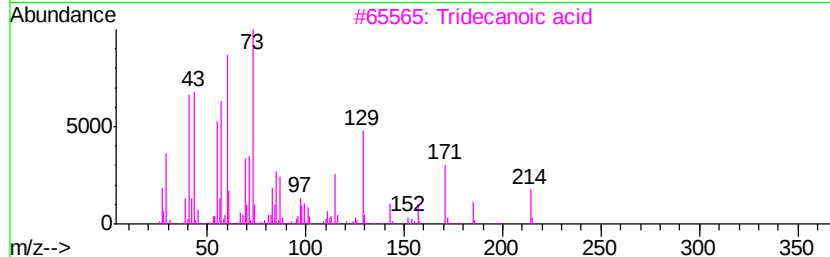
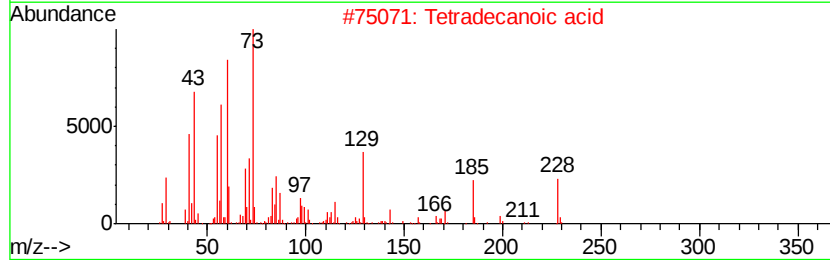
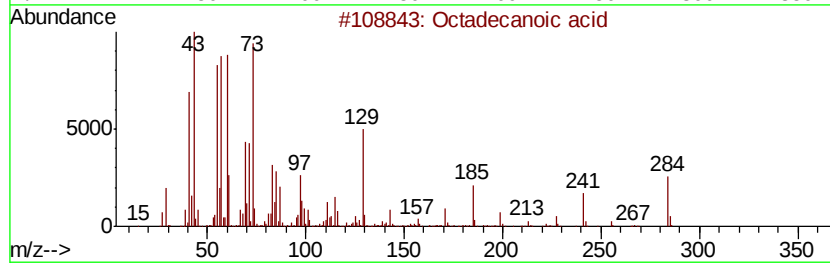
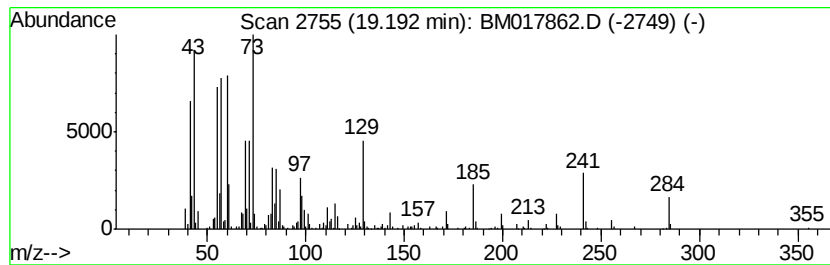
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Peak Number 5 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.19	8.19 ng	1081630	Chrysene-d12	21.19

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3		Tridecanoic acid	214	C13H26O2	000638-53-9	74
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	70
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64



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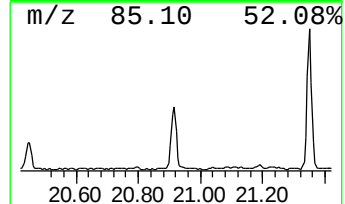
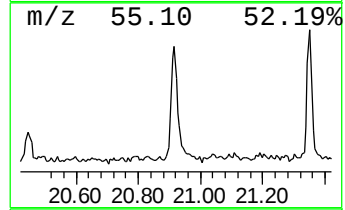
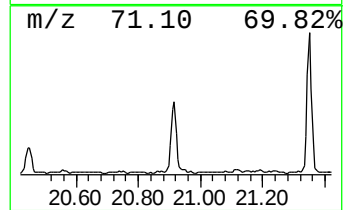
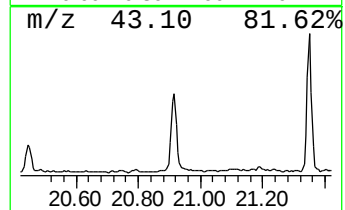
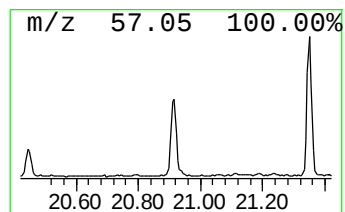
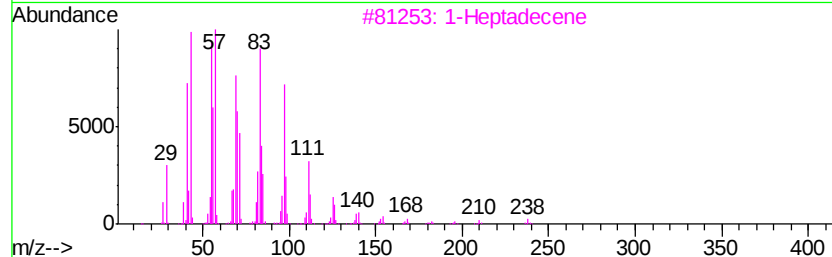
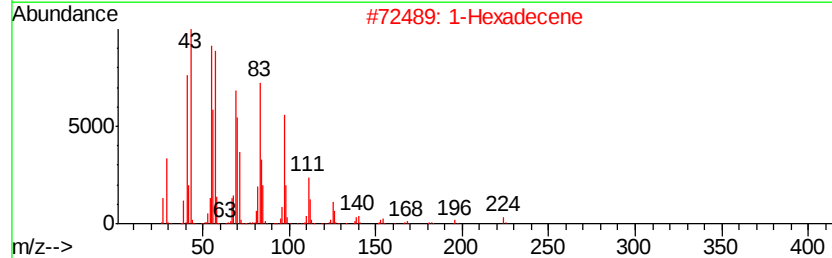
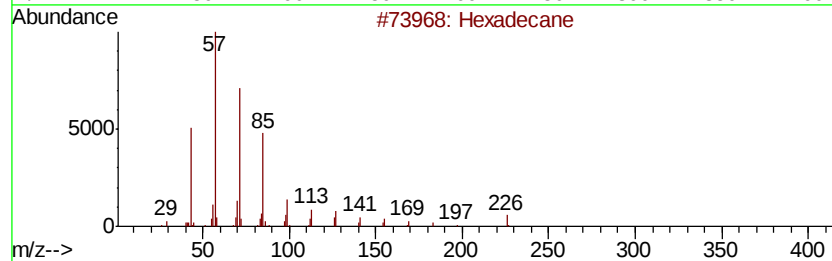
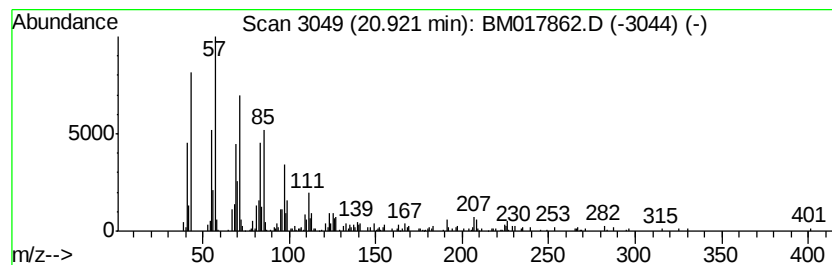
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	4.15 ng	547364	Chrysene-d12	21.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	96
2		1-Hexadecene	224	C16H32	000629-73-2	94
3		1-Heptadecene	238	C17H34	006765-39-5	94
4		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	91
5		Tritetracontane	605	C43H88	007098-21-7	87



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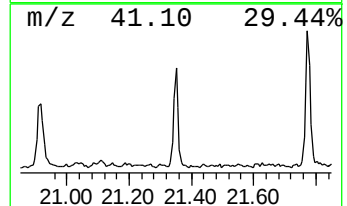
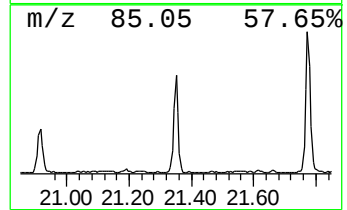
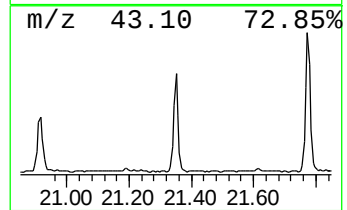
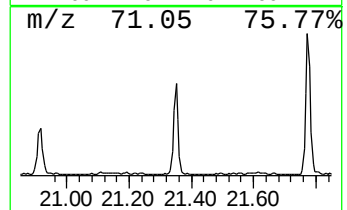
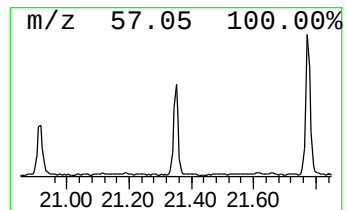
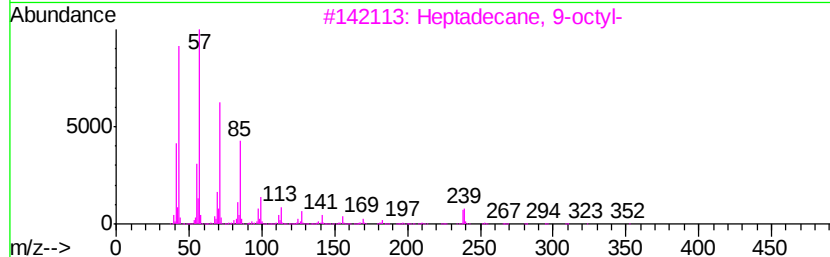
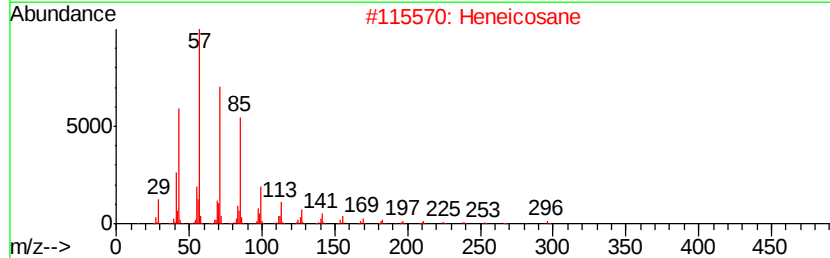
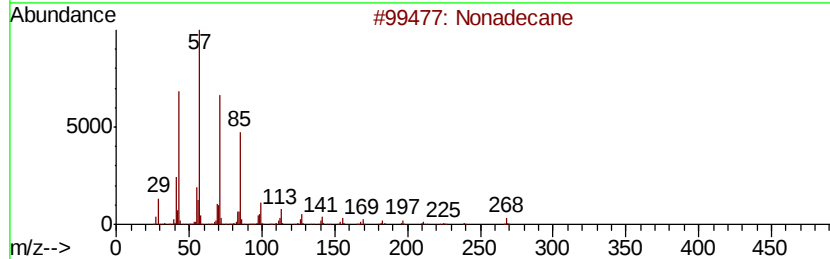
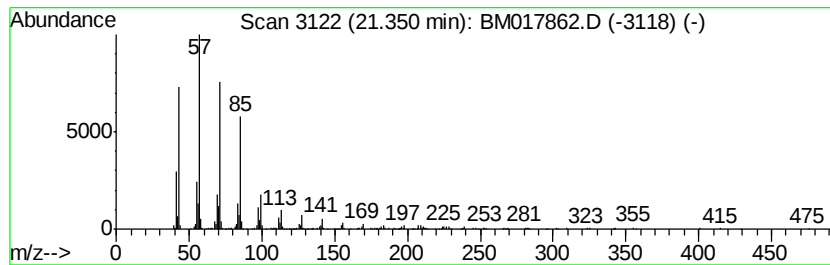
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	4.57 ng	603518	Chrysene-d12	21.19
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonadecane	268 C19H40	000629-92-5	94
2	Heneicosane	296 C21H44	000629-94-7	91
3	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
4	Pentadecane, 8-heptyl-	310 C22H46	071005-15-7	91
5	Tetracosane	338 C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
Data File : BM017862.D
Acq On : 01 Dec 2018 18:23
Operator : JU/SJ
Sample : J6060-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

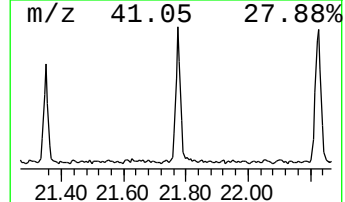
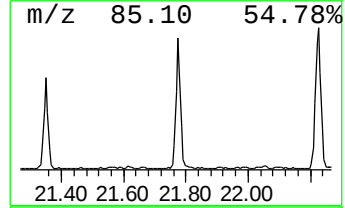
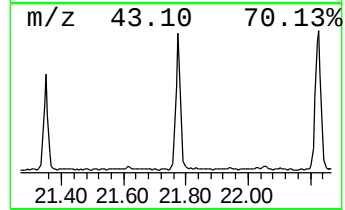
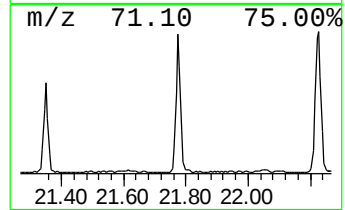
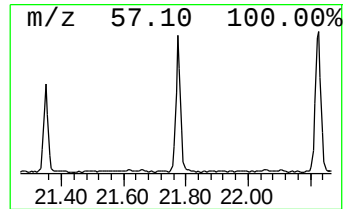
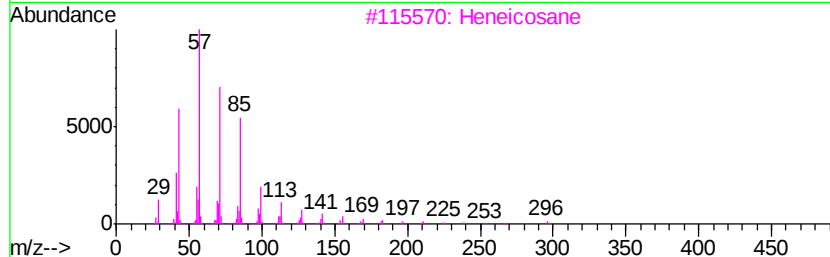
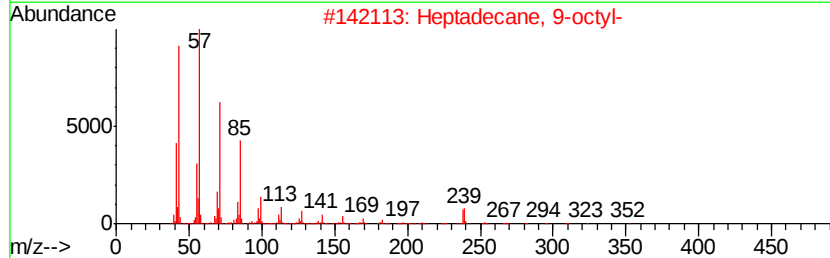
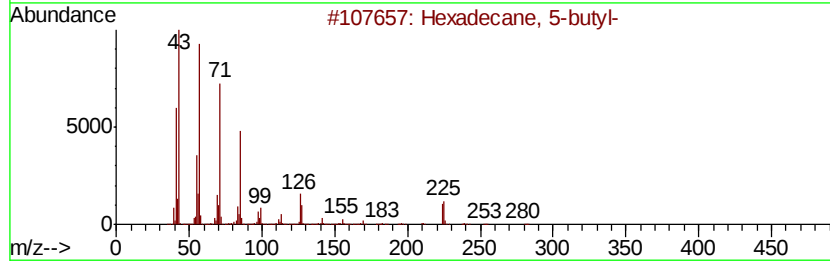
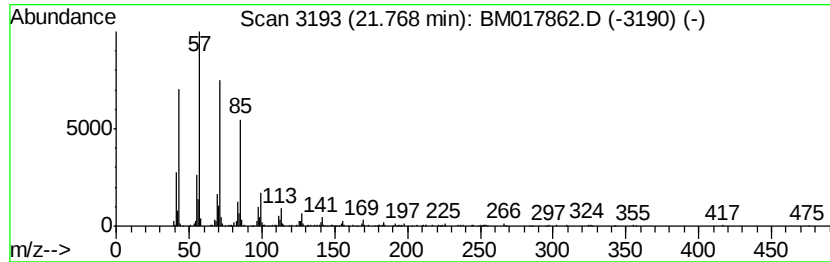
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane, 5-butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.77	6.88 ng	908201	Chrysene-d12		21.19
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 5-butyl-	282	C20H42	006912-07-8	93
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Tetratetracontane	619	C44H90	007098-22-8	91
5	Heptadecane	240	C17H36	000629-78-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
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Acq On : 01 Dec 2018 18:23
Operator : JU/SJ
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Instrument :
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ClientSampleId :
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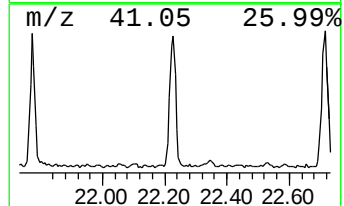
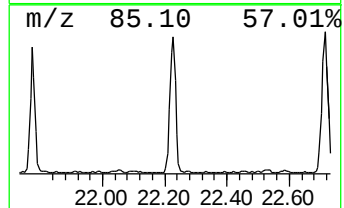
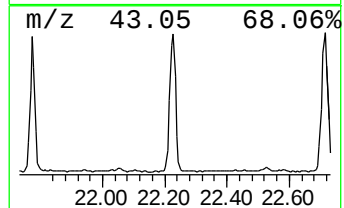
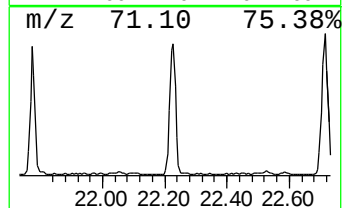
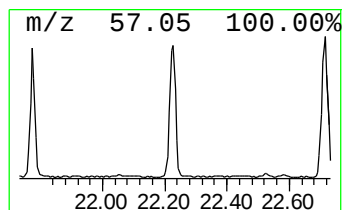
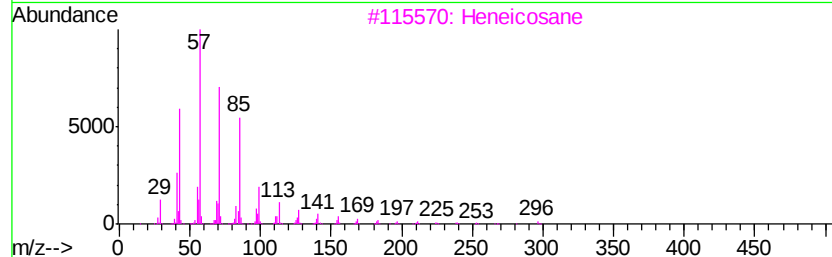
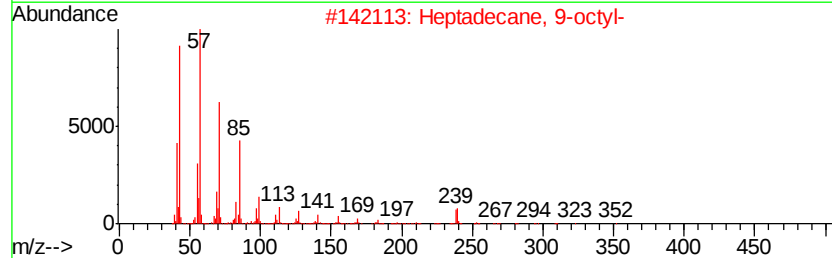
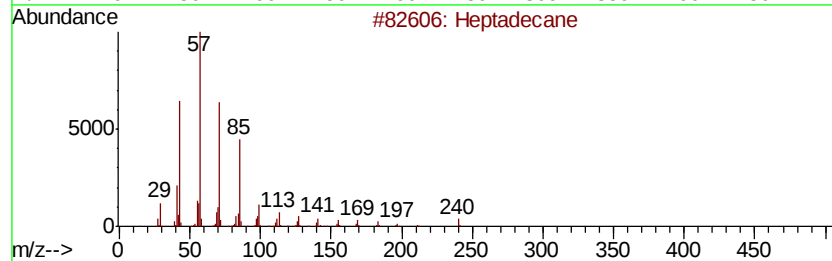
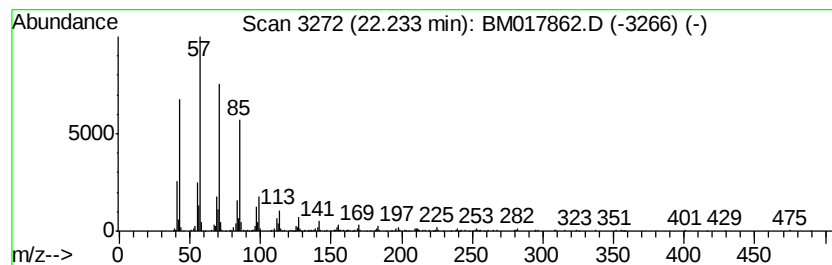
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.23	8.70 ng	1149050	Chrysene-d12	21.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	94
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Eicosane	282	C20H42	000112-95-8	91
5		Nonacosane	408	C29H60	000630-03-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
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ClientSampleId :
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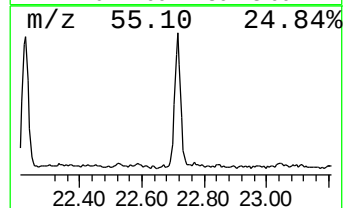
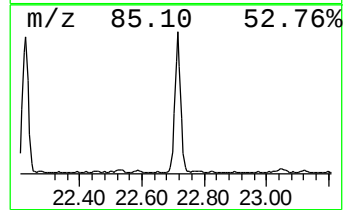
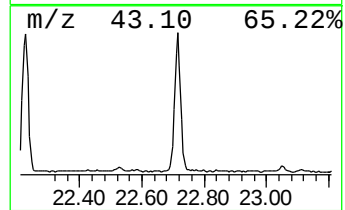
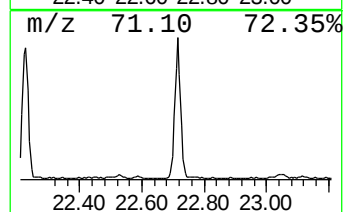
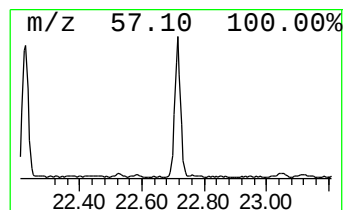
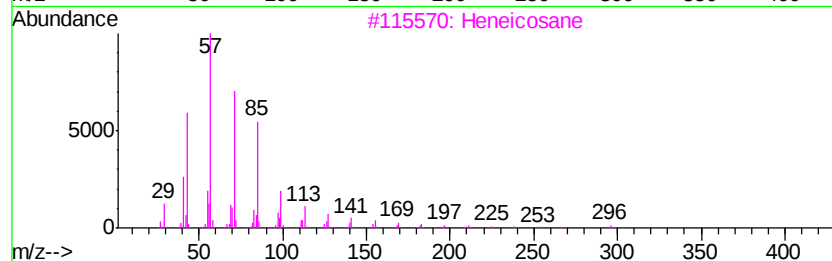
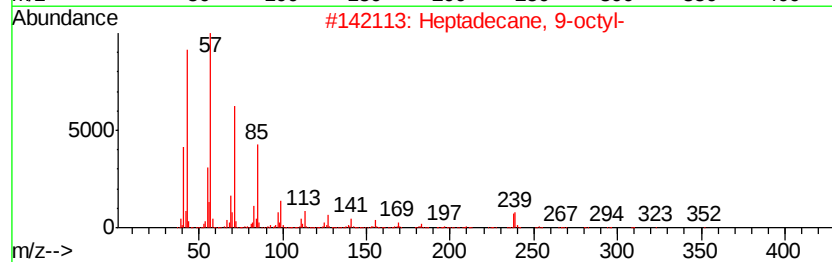
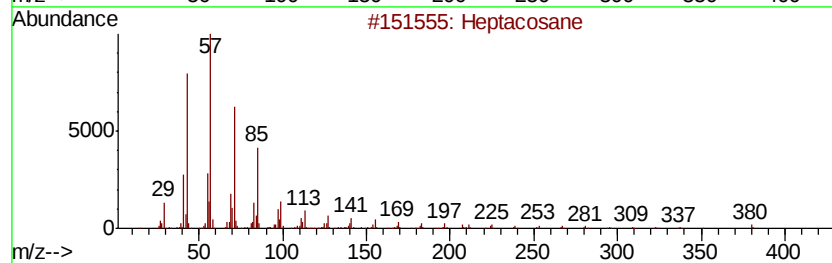
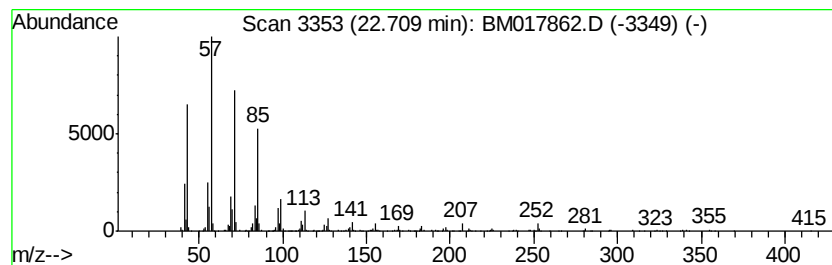
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.71	10.17 ng	1322350	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Nonacosane	408	C29H60	000630-03-5	90
5		Docosane	310	C22H46	000629-97-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
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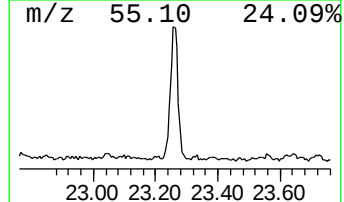
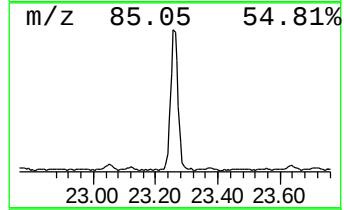
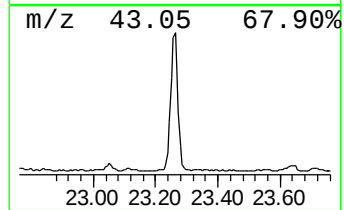
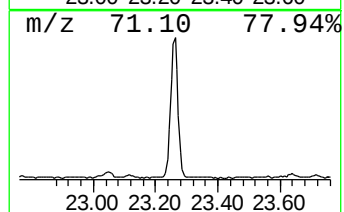
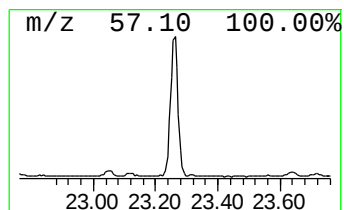
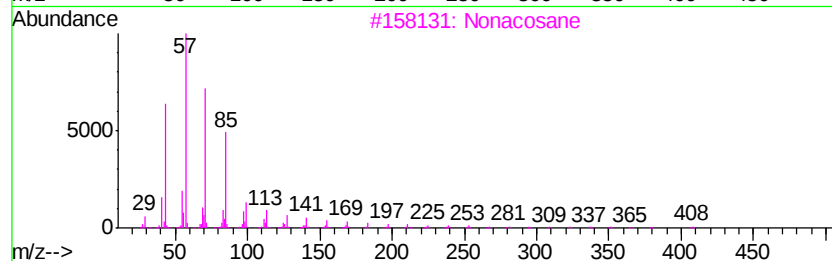
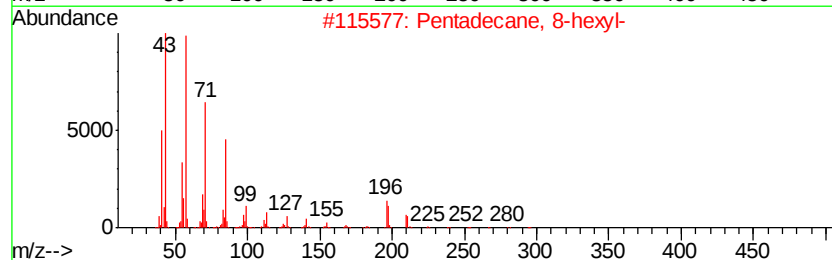
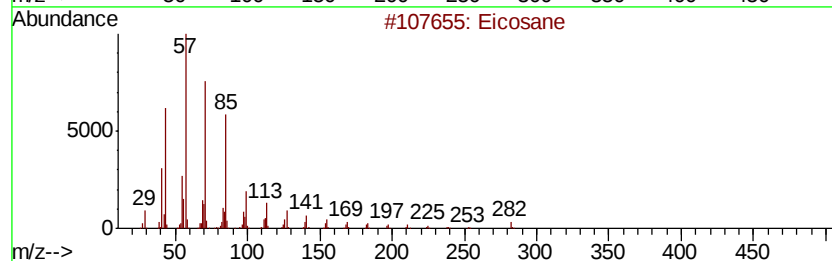
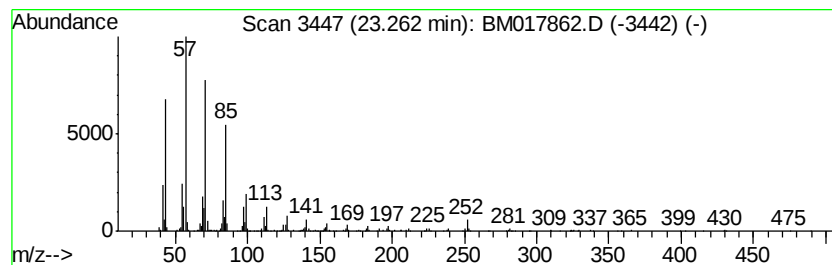
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.26	9.36 ng	1217700	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	90
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
3		Nonacosane	408	C29H60	000630-03-5	90
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
5		Tetratetracontane	619	C44H90	007098-22-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
Data File : BM017862.D
Acq On : 01 Dec 2018 18:23
Operator : JU/SJ
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20180947-COMPOSITE-6

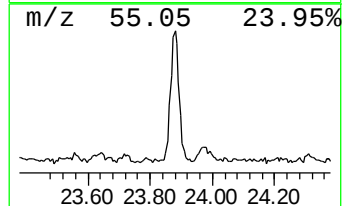
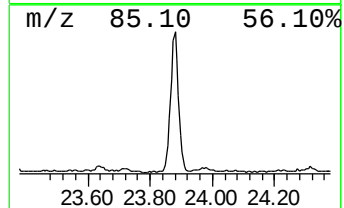
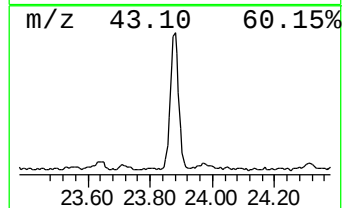
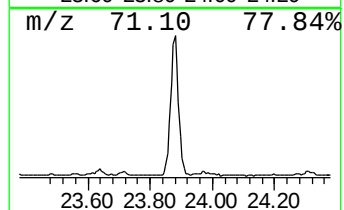
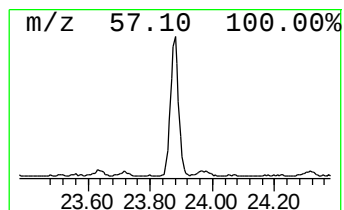
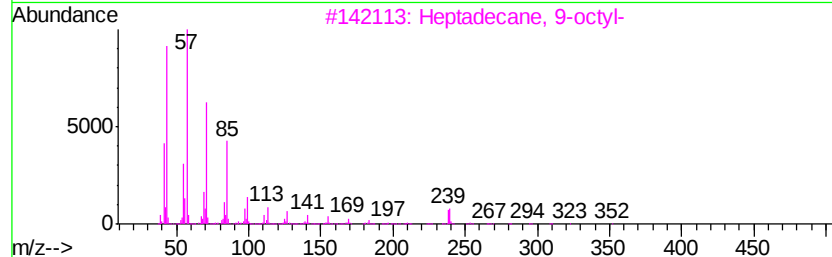
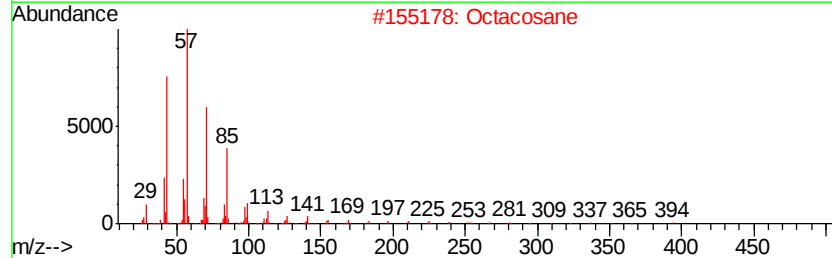
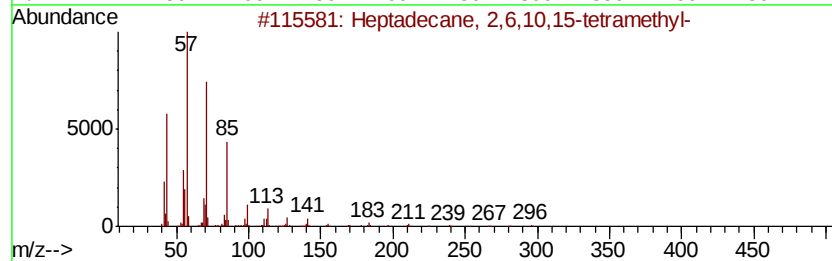
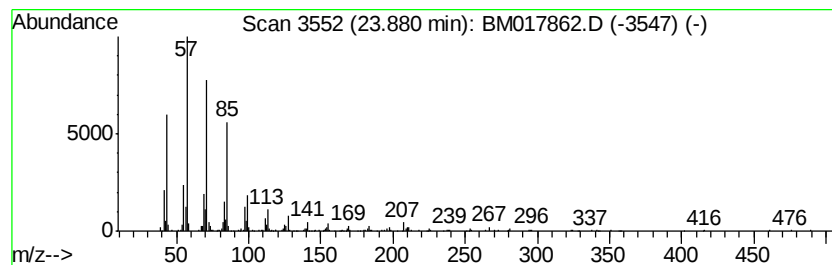
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane, 2,6,10,15-tetr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.88	7.35 ng	955456	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2		Octacosane	394	C28H58	000630-02-4	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Hexadecane	226	C16H34	000544-76-3	91
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91



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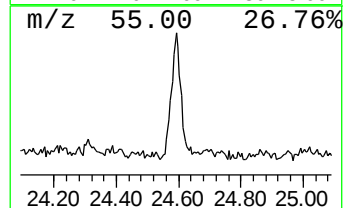
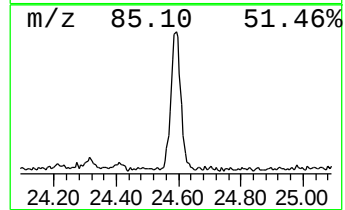
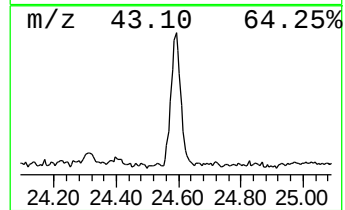
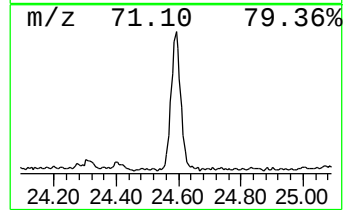
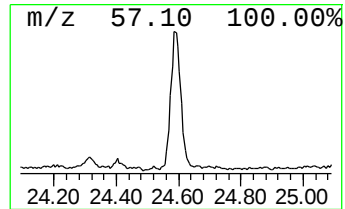
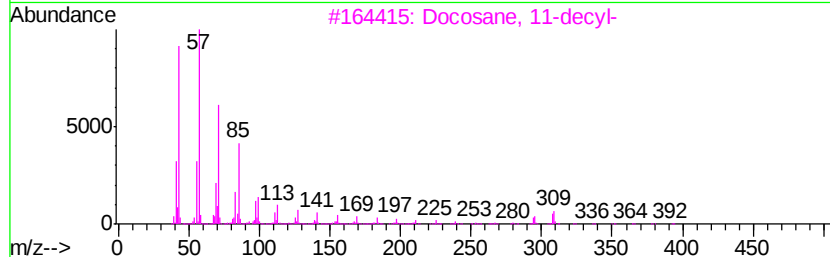
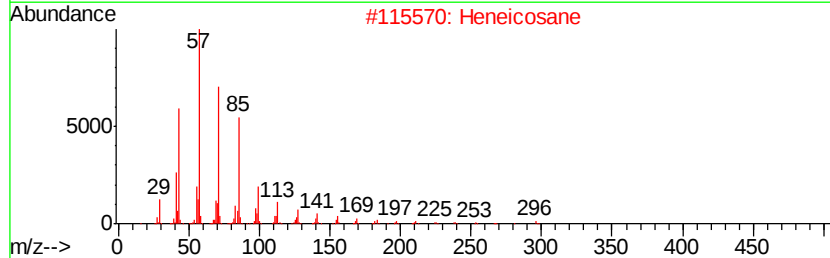
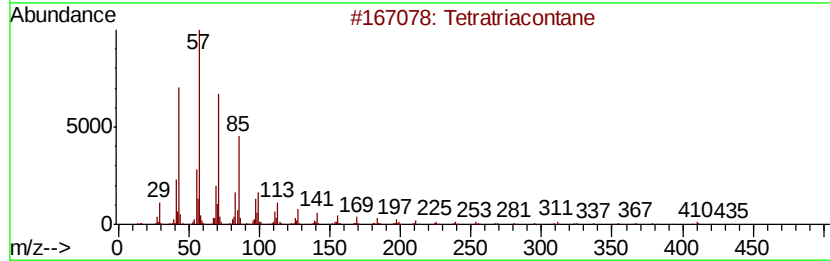
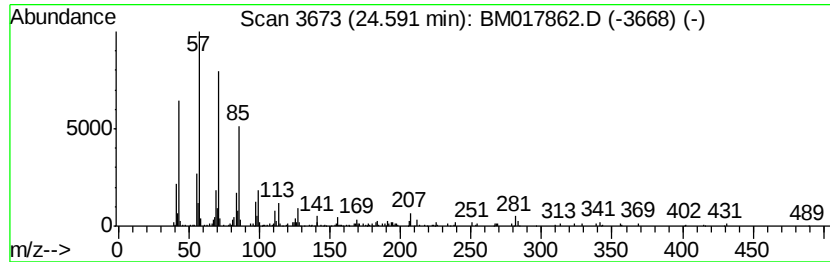
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetratriacontane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
24.59	3.81 ng	495056	Perylene-d12		23.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratriacontane	479	C34H70	014167-59-0	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Docosane, 11-decyl-	451	C32H66	055401-55-3	91
4	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM120218\
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Sample : J6060-06
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20180947-COMPOSITE-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM111218.M
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
2-Pentanone, 4-hy...	4.77	3.3	ng	254436	1	7.58	1542930	20.0
unknown7.12	7.12	102.9	ng	7939800	1	7.58	1542930	20.0
n-Hexadecanoic acid	17.90	10.5	ng	1623640	4	16.99	3088020	20.0
Octadecanoic acid	19.19	8.2	ng	1081630	5	21.19	2640770	20.0
Hexadecane	20.92	4.2	ng	547364	5	21.19	2640770	20.0
Nonadecane	21.35	4.6	ng	603518	5	21.19	2640770	20.0
Hexadecane, 5-butyl-	21.77	6.9	ng	908201	5	21.19	2640770	20.0
Heptadecane	22.23	8.7	ng	1149050	5	21.19	2640770	20.0
Heptacosane	22.71	10.2	ng	1322350	6	23.37	2601480	20.0
Eicosane	23.26	9.4	ng	1217700	6	23.37	2601480	20.0
Heptadecane, 2,6,...	23.88	7.3	ng	955456	6	23.37	2601480	20.0
Tetratriacontane	24.59	3.8	ng	495056	6	23.37	2601480	20.0