

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
 Data File : BM018613.D
 Acq On : 17 Jan 2019 14:36
 Operator : JU/SJ
 Sample : K1147-22
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-8

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_M\METHODS\8270-BM010919.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	3.516	70	73	87	rBV	226338	344624	1.93%	0.233%
2	4.887	300	306	314	rBV	326018	504902	2.83%	0.341%
3	5.334	375	382	394	rBV	7417346	10405200	58.36%	7.021%
4	6.922	644	652	666	rBV	6642095	11256666	63.13%	7.596%
5	7.275	703	712	722	rBV	7492886	12070398	67.69%	8.145%
6	7.739	785	791	798	rVB	1452513	2176123	12.20%	1.468%
7	8.057	837	845	855	rVB	6085223	9584496	53.75%	6.468%
8	8.910	981	990	1003	rBV	4344181	7101820	39.83%	4.792%
9	10.528	1258	1265	1273	rBV	1652337	2783922	15.61%	1.879%
10	13.004	1679	1686	1701	rBV	10556139	15730015	88.22%	10.615%
11	13.845	1822	1829	1841	rBV	878376	1212912	6.80%	0.818%
12	14.386	1915	1921	1930	rVB2	2178131	3363567	18.86%	2.270%
13	15.886	2168	2176	2193	rBV	7914409	11302143	63.39%	7.627%
14	17.139	2382	2389	2395	rBV	2525612	3609194	20.24%	2.435%
15	18.027	2534	2540	2558	rBV2	1188341	2042462	11.45%	1.378%
16	19.209	2736	2741	2746	rVB3	129771	191036	1.07%	0.129%
17	19.309	2753	2758	2773	rBV	680174	1227620	6.88%	0.828%
18	19.774	2831	2837	2847	rBV	14487745	17830676	100.00%	12.032%
19	21.033	3047	3051	3057	rBV	569059	675941	3.79%	0.456%
20	21.333	3096	3102	3106	rBV	2793525	3844639	21.56%	2.594%
21	21.468	3121	3125	3132	rVB	742359	857920	4.81%	0.579%
22	21.909	3195	3200	3207	rBV	1993279	2390351	13.41%	1.613%
23	22.374	3274	3279	3295	rVB	3417450	4776093	26.79%	3.223%
24	22.891	3361	3367	3376	rBV	3981416	6079172	34.09%	4.102%
25	23.468	3459	3465	3477	rVB	3564919	6150499	34.49%	4.150%
26	23.603	3482	3488	3499	rVB	1894249	3723429	20.88%	2.513%
27	24.121	3570	3576	3586	rBV	2335432	4575137	25.66%	3.087%
28	24.885	3699	3706	3715	rBV	1083052	2380694	13.35%	1.606%

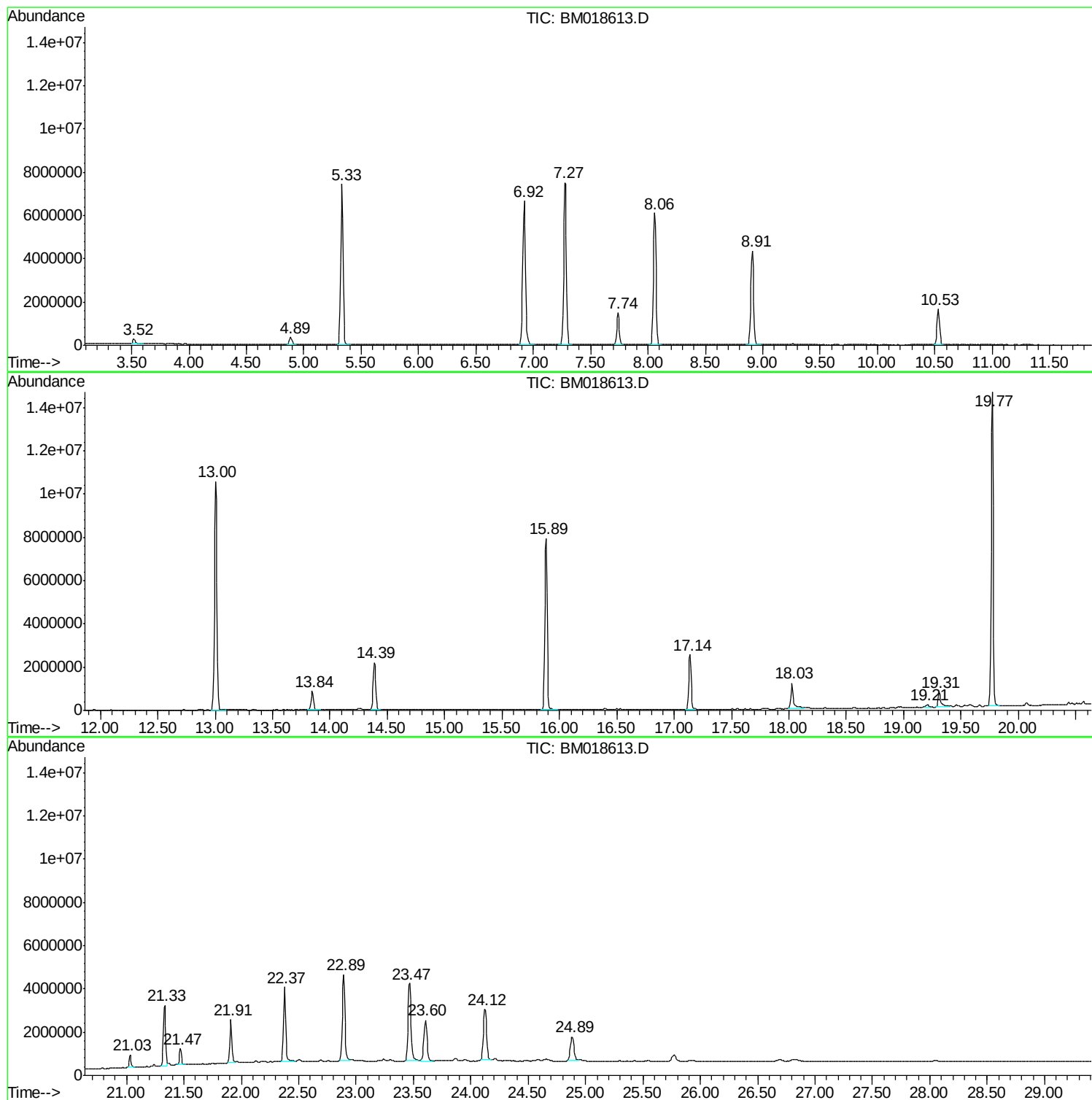
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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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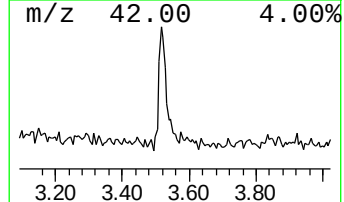
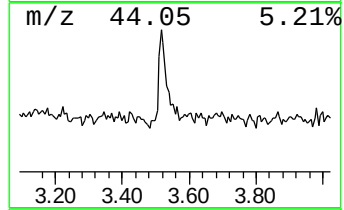
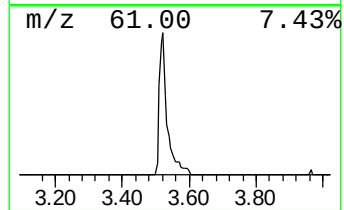
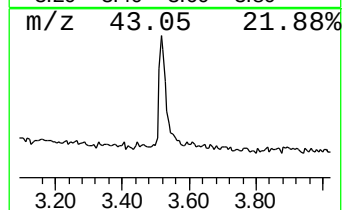
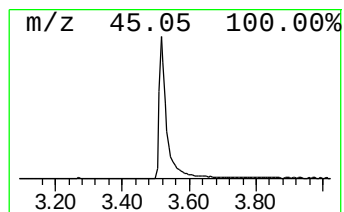
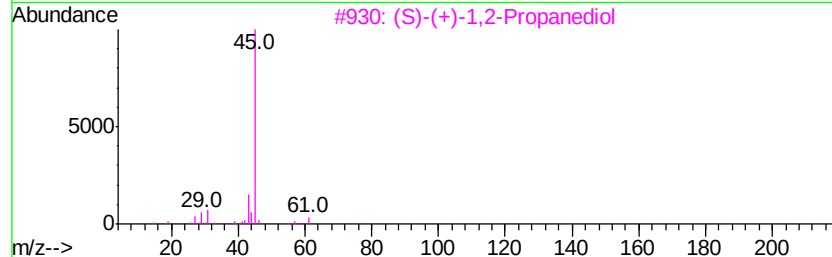
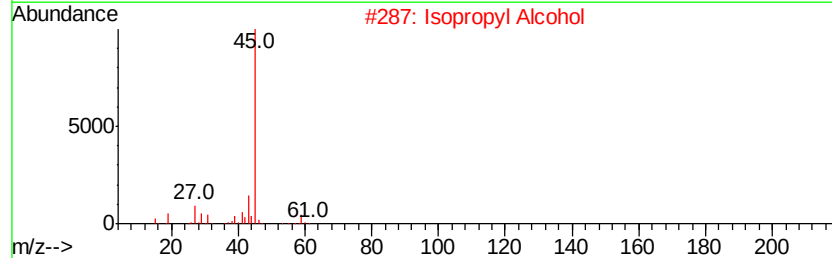
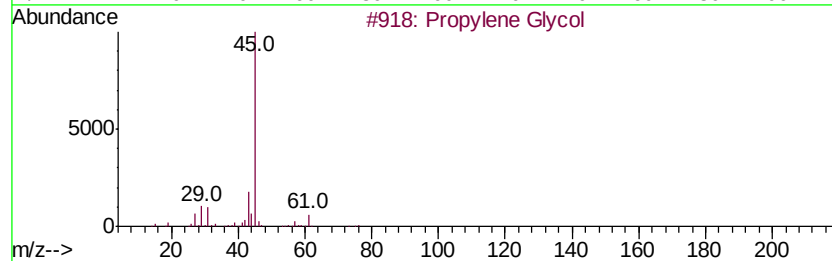
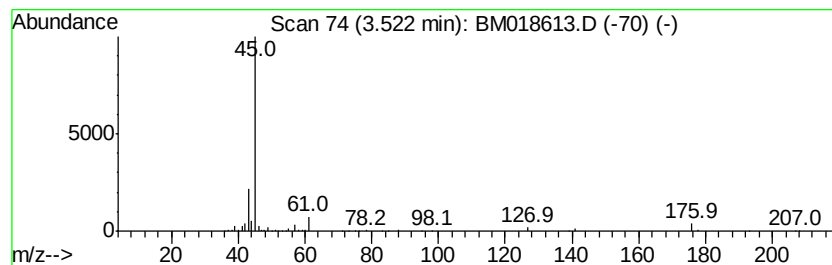
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propylene Glycol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	3.17 ng	344624	1,4-Dichlorobenzene-d4	7.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propylene Glycol	76	C3H8O2	000057-55-6	86
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
4		2-Heptanol, (S)-	116	C7H16O	006033-23-4	9
5		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9



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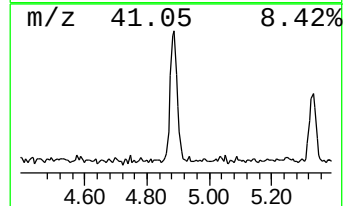
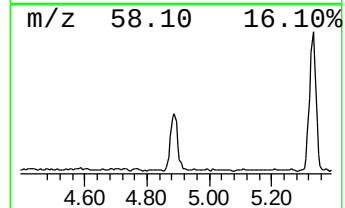
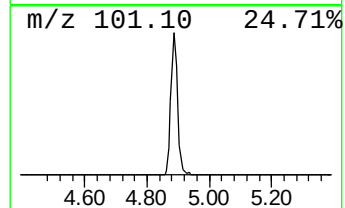
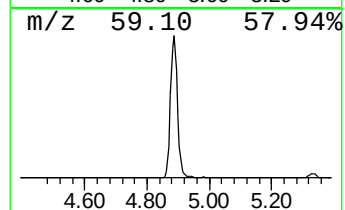
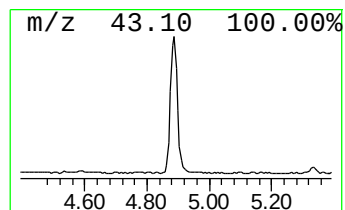
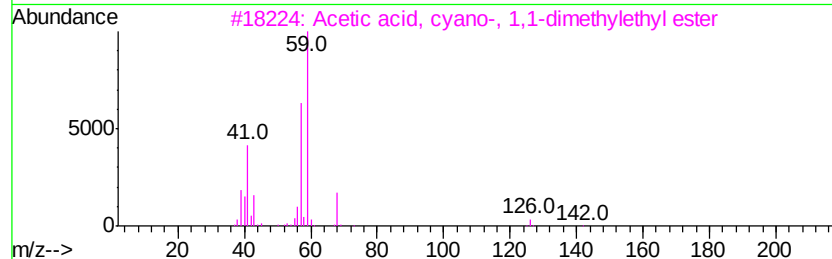
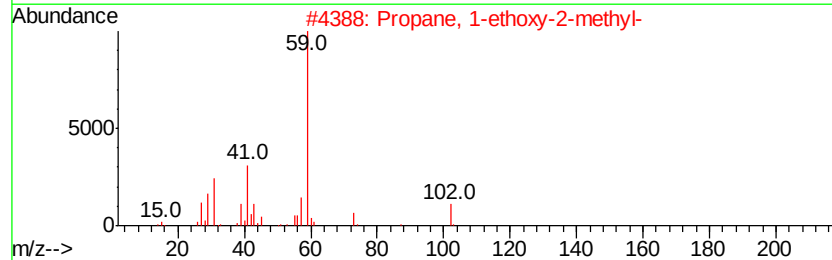
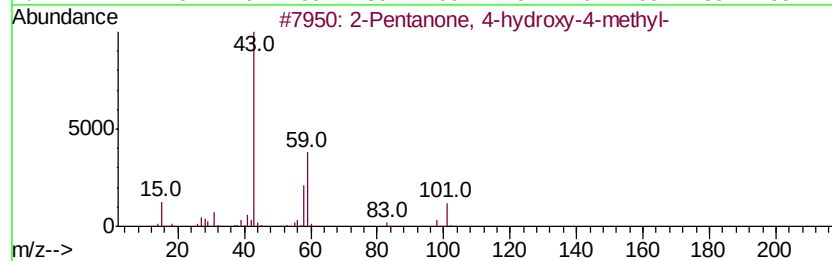
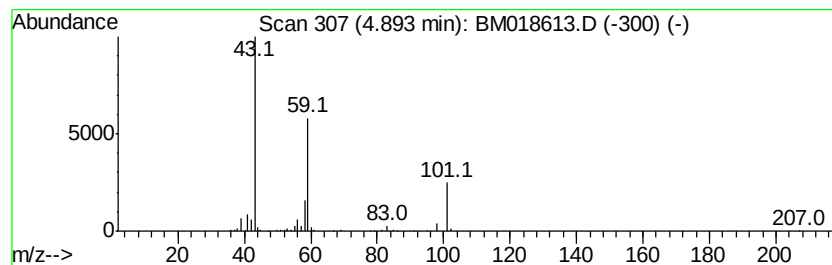
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	4.64 ng	504902	1,4-Dichlorobenzene-d4	7.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	38
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
4			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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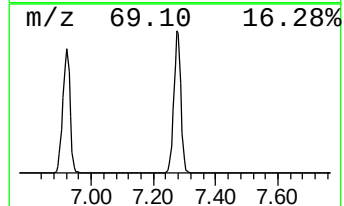
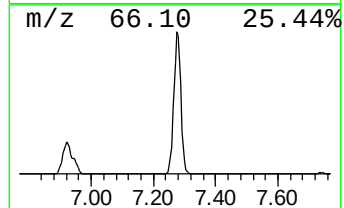
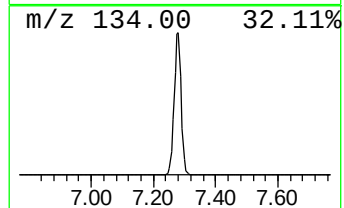
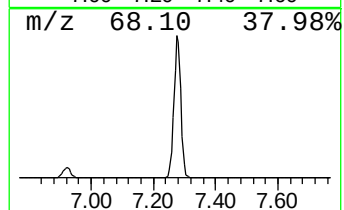
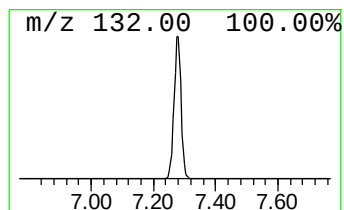
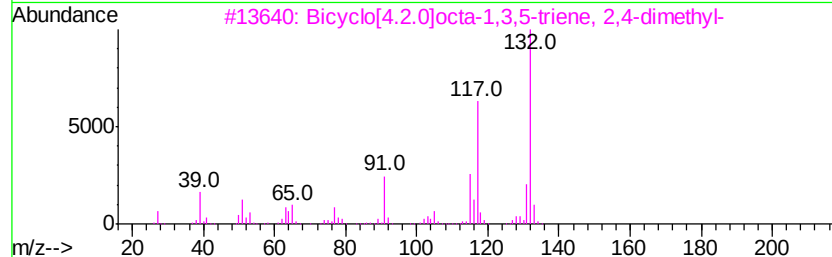
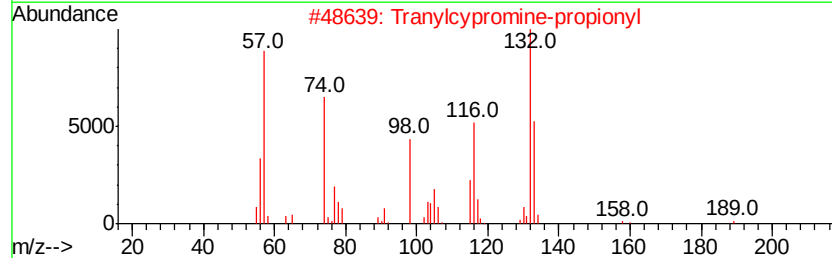
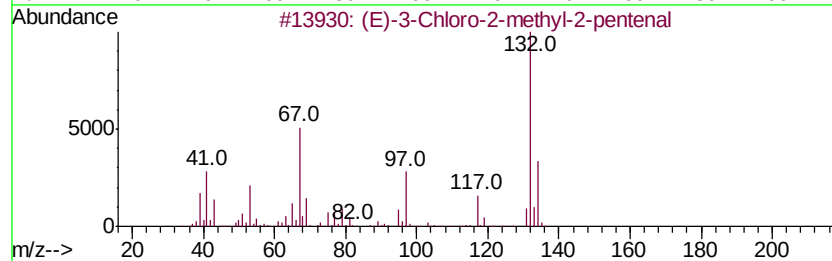
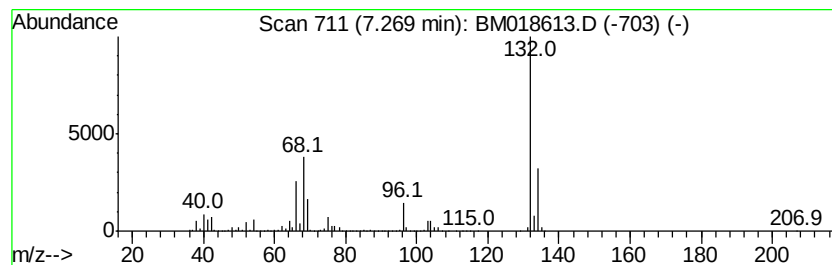
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Peak Number 3 unknown7.27 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	110.94 ng	12070400	1,4-Dichlorobenzene-d4	7.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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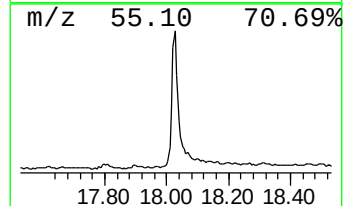
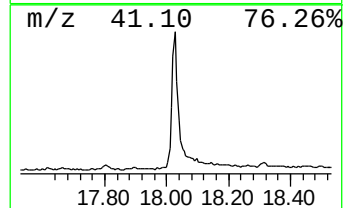
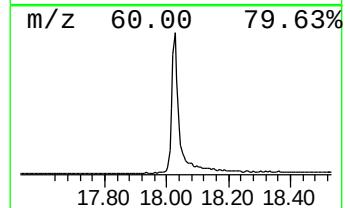
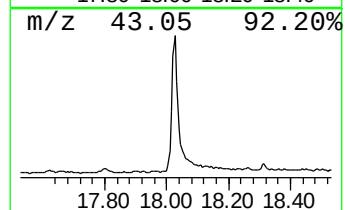
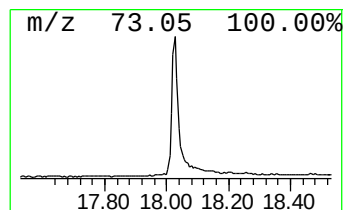
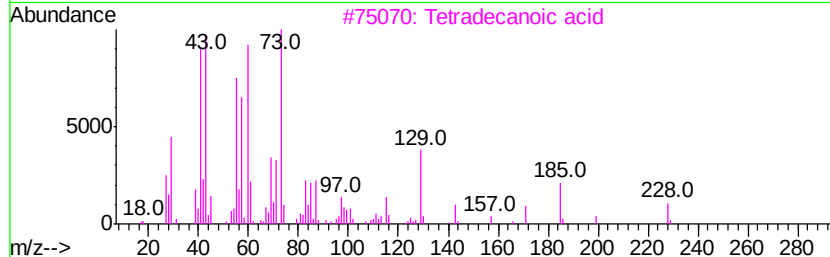
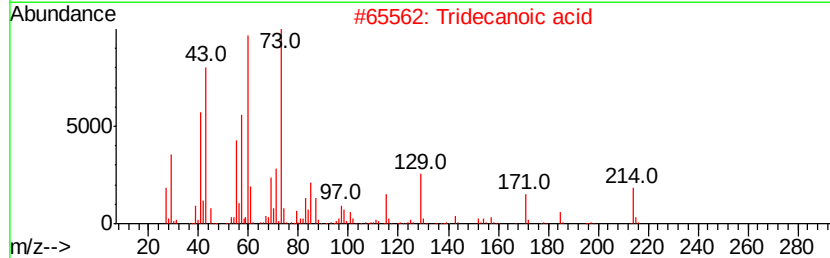
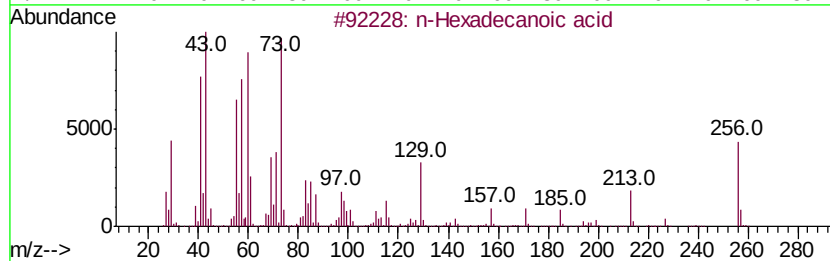
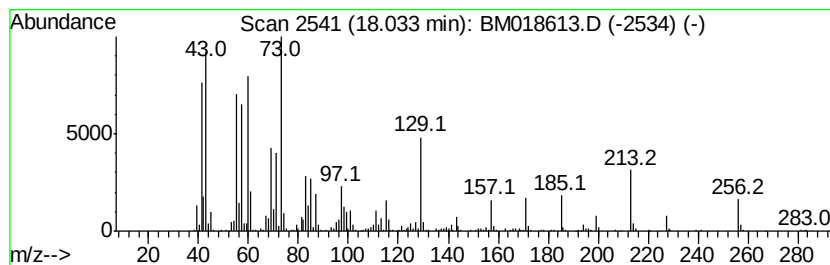
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	11.32 ng	2042460	Phenanthrene-d10	17.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4		Undecanoic acid	186	C11H22O2	000112-37-8	83
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	83



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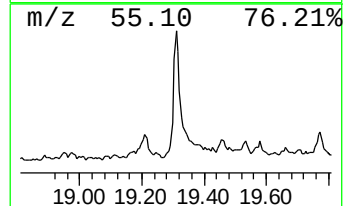
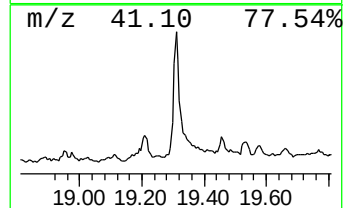
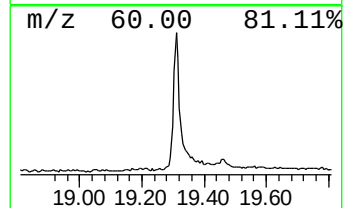
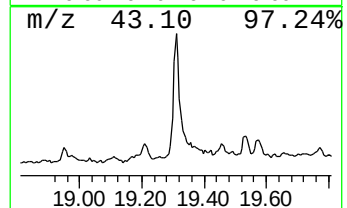
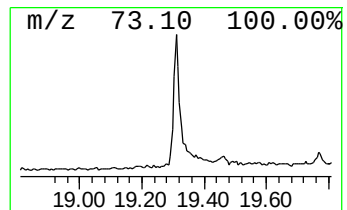
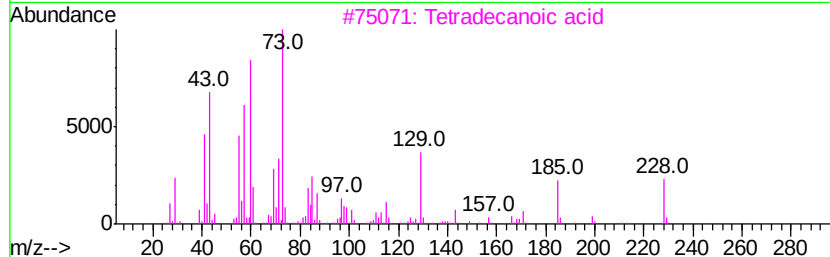
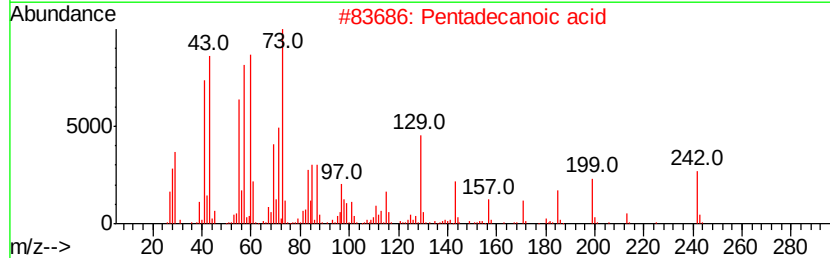
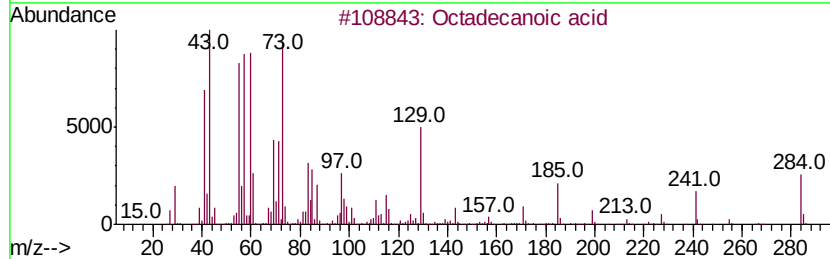
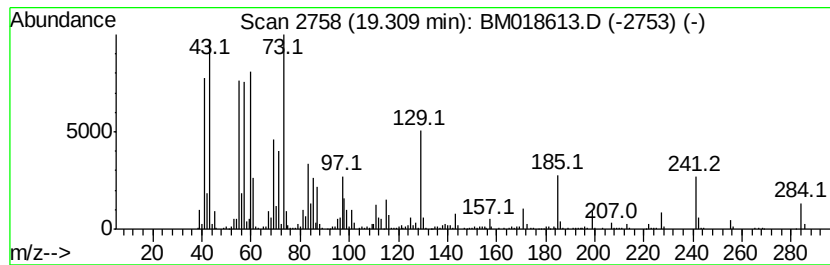
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Peak Number 6 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.31	6.39 ng	1227620	Chrysene-d12	21.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	89
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4	Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	70



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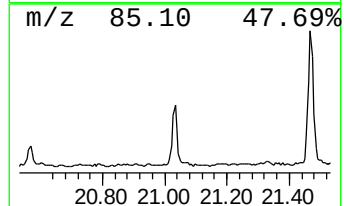
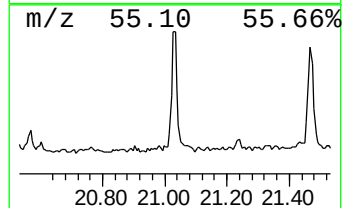
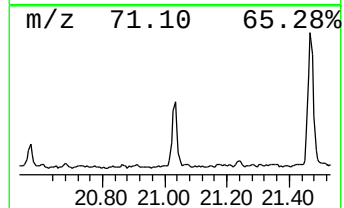
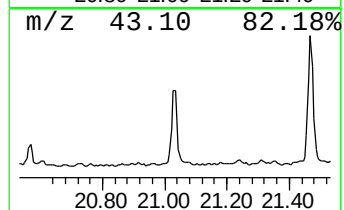
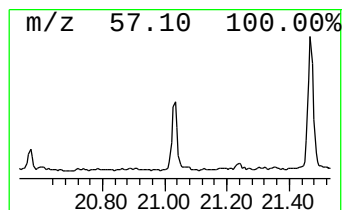
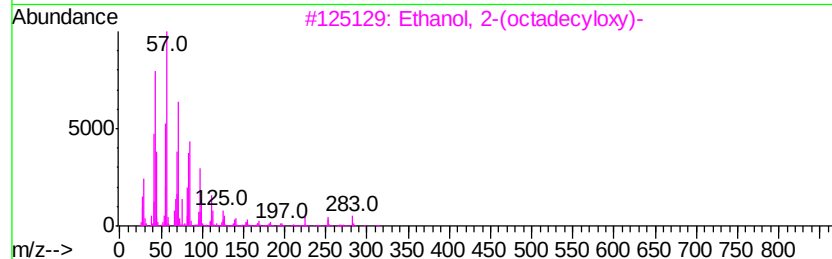
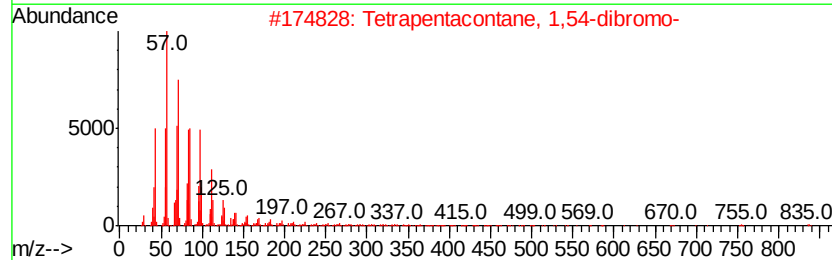
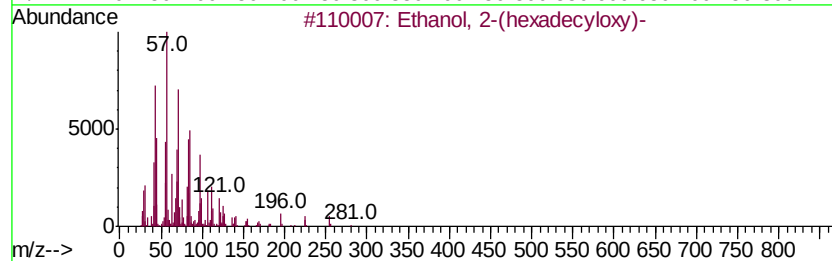
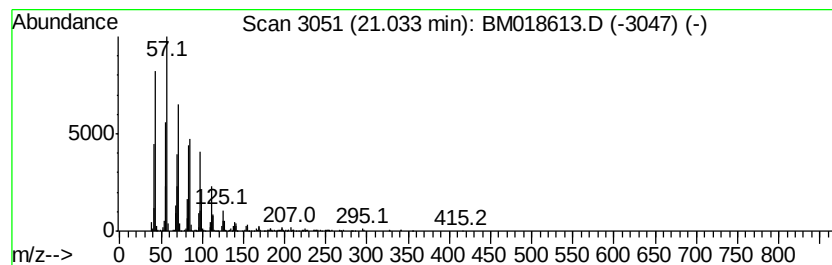
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Ethanol, 2-(hexadecyloxy)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.03	3.52 ng	675941	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	91
2		Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	90
3		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	87
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87
5		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
Data File : BM018613.D
Acq On : 17 Jan 2019 14:36
Operator : JU/SJ
Sample : K1147-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
TP-8

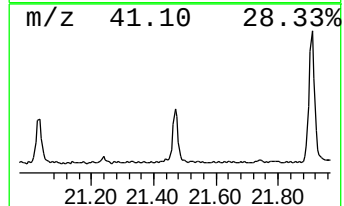
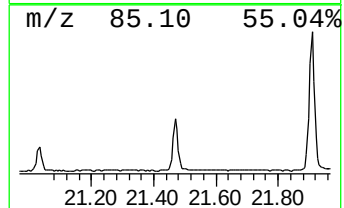
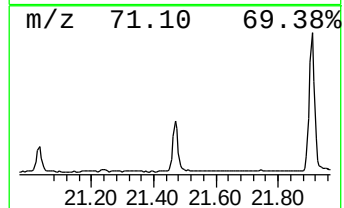
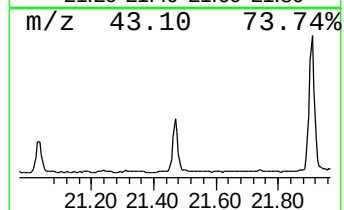
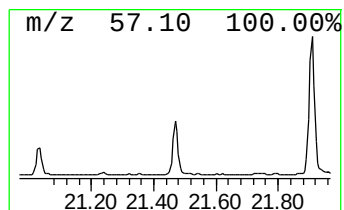
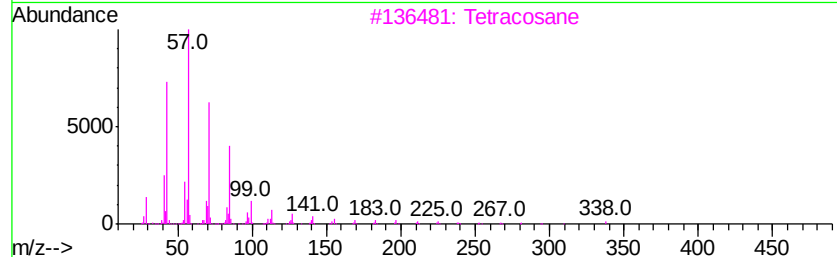
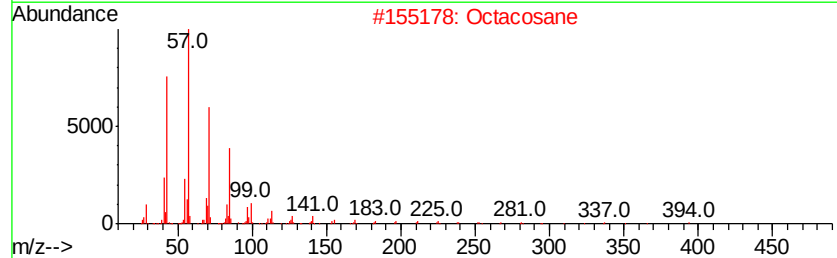
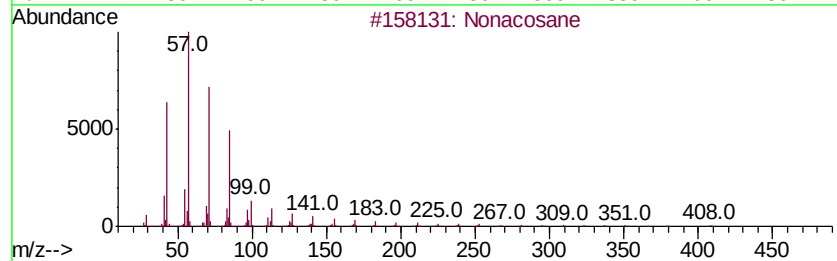
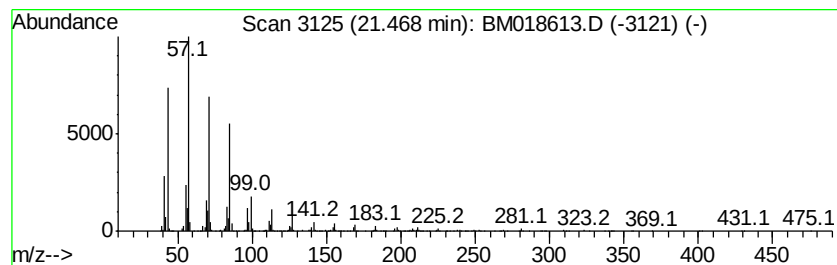
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.47	4.46 ng	857920	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonacosane	408	C29H60	000630-03-5	90
2		Octacosane	394	C28H58	000630-02-4	90
3		Tetracosane	338	C24H50	000646-31-1	90
4		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
Data File : BM018613.D
Acq On : 17 Jan 2019 14:36
Operator : JU/SJ
Sample : K1147-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

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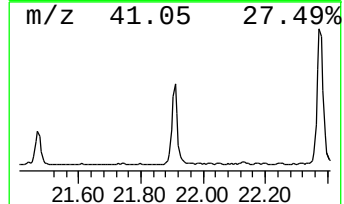
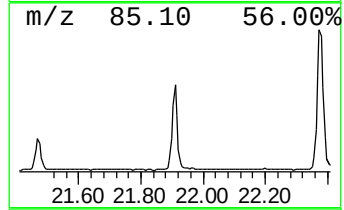
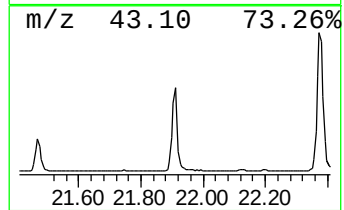
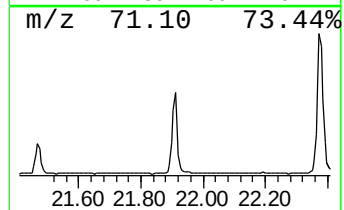
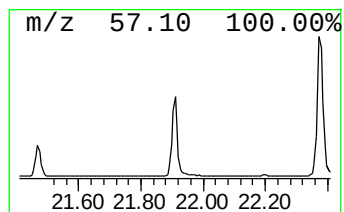
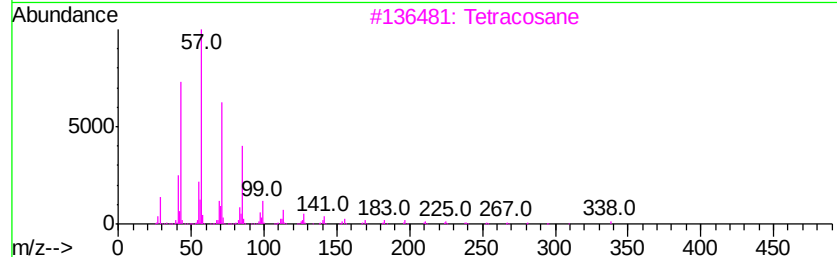
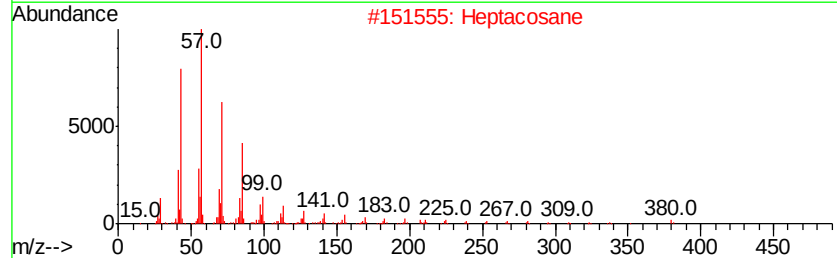
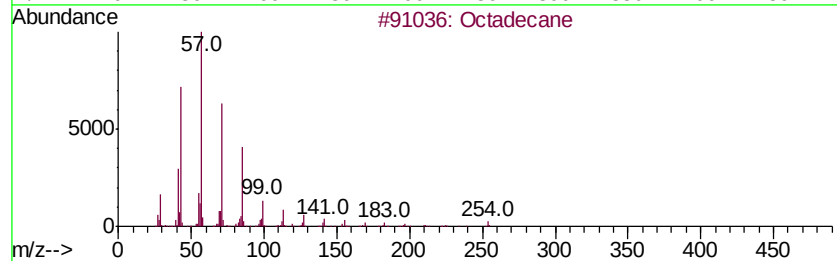
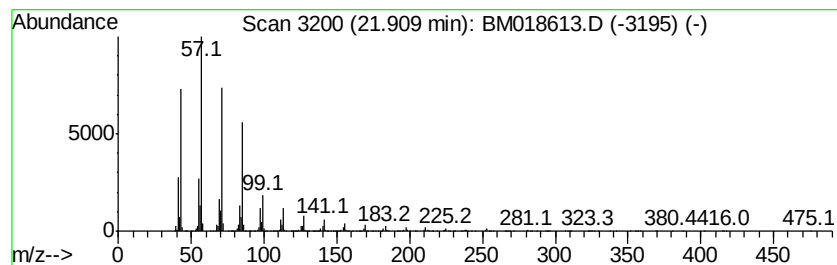
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.91	12.43 ng	2390350	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	97
2		Heptacosane	380	C27H56	000593-49-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
 Data File : BM018613.D
 Acq On : 17 Jan 2019 14:36
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 Sample : K1147-22
 Misc :
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Instrument :
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 ClientSampleId :
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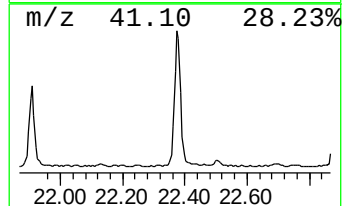
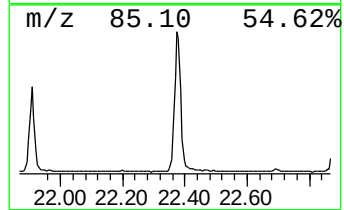
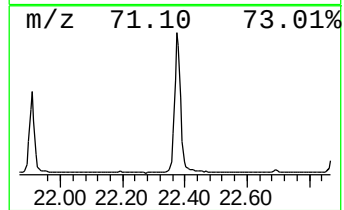
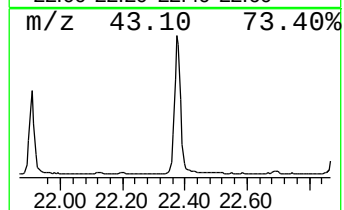
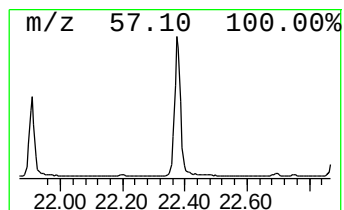
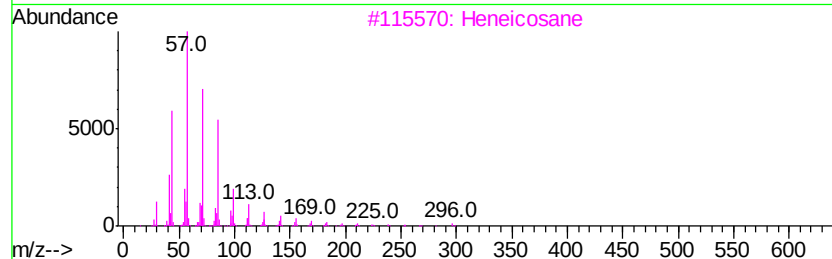
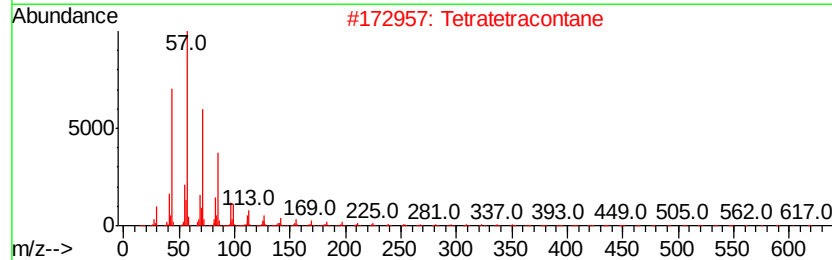
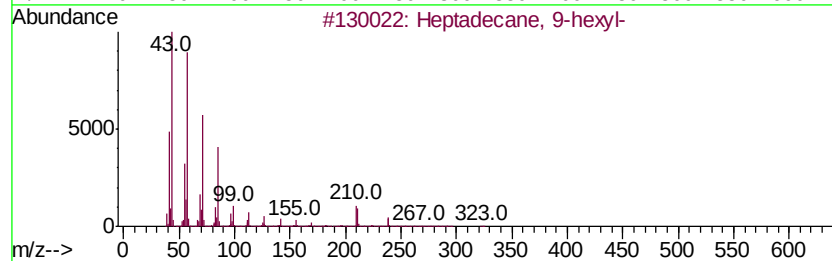
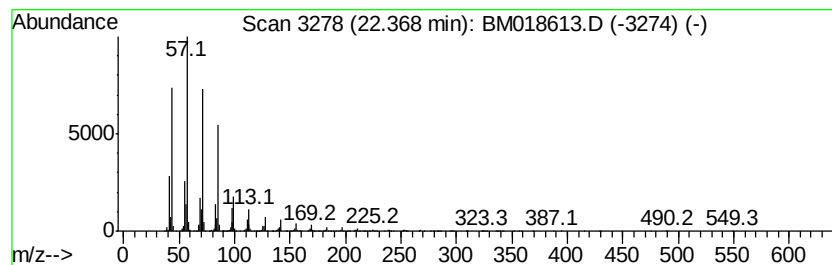
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 Heptadecane, 9-hexyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.37	24.85 ng	4776090	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Tetratriacontane	479	C34H70	014167-59-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
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Sample : K1147-22
Misc :
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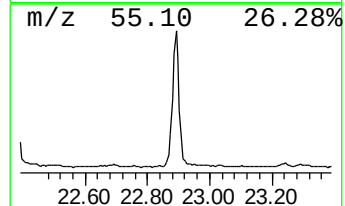
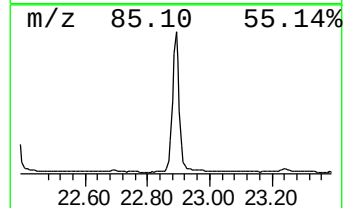
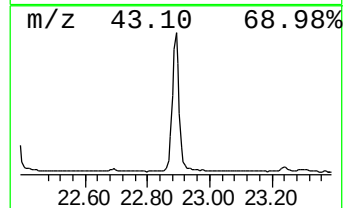
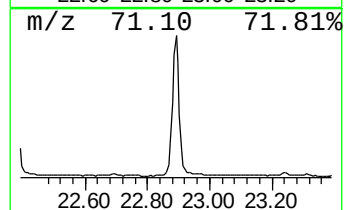
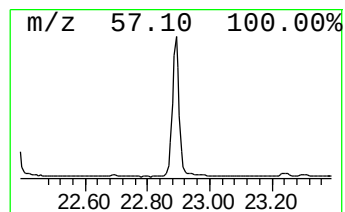
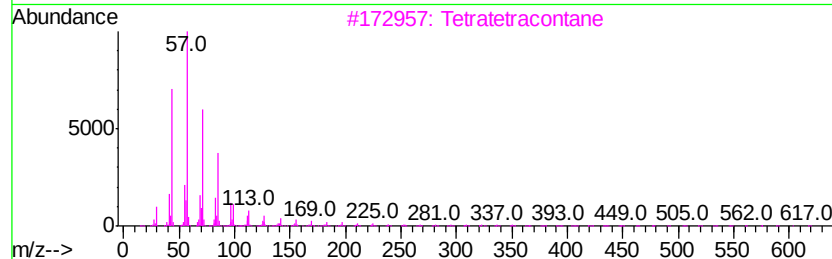
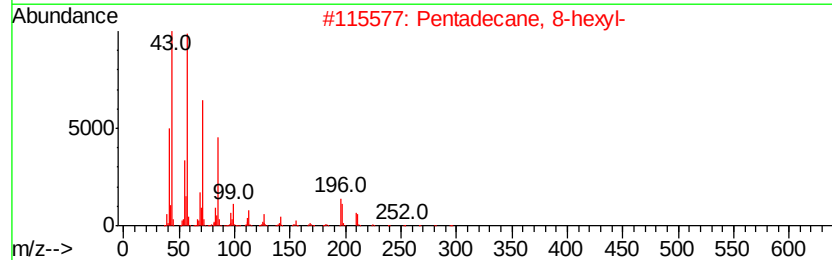
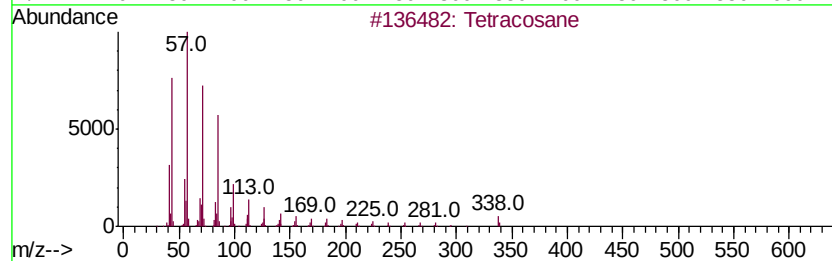
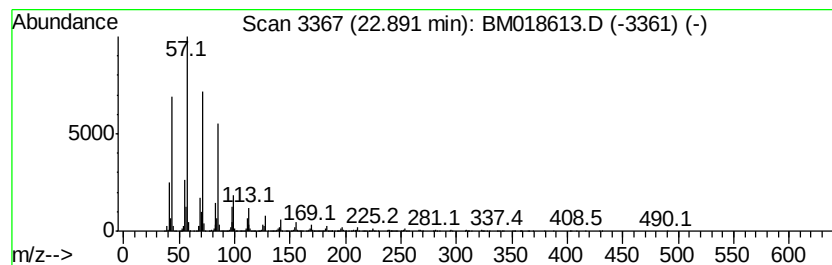
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.89	32.65 ng	6079170	Perylene-d12	23.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetracosane	338 C24H50	000646-31-1 98
2		Pentadecane, 8-hexyl-	296 C21H44	013475-75-7 94
3		Tetratetracontane	619 C44H90	007098-22-8 91
4		Nonacosane	408 C29H60	000630-03-5 91
5		Heptadecane, 9-octyl-	352 C25H52	007225-64-1 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
Data File : BM018613.D
Acq On : 17 Jan 2019 14:36
Operator : JU/SJ
Sample : K1147-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

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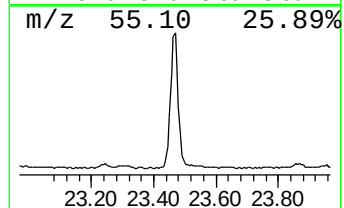
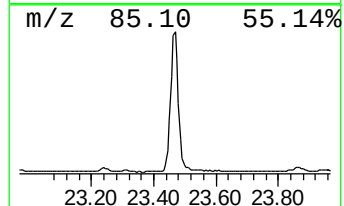
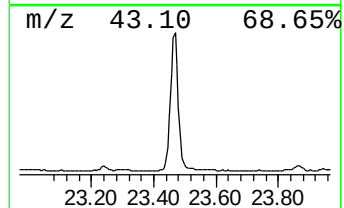
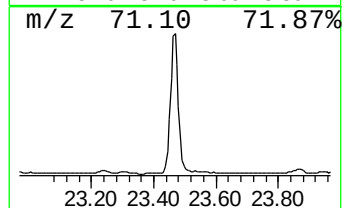
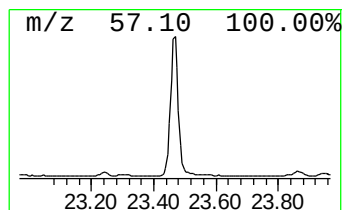
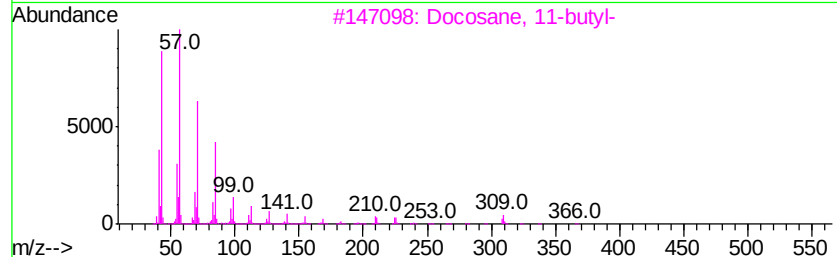
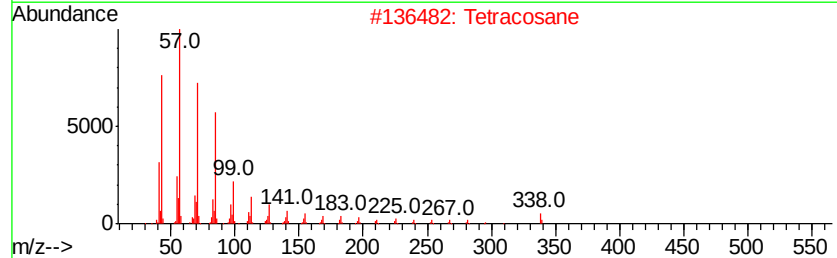
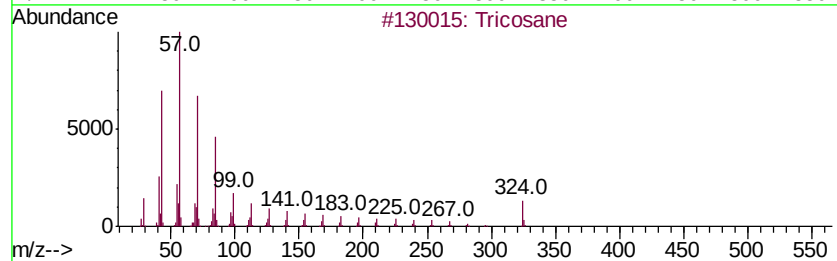
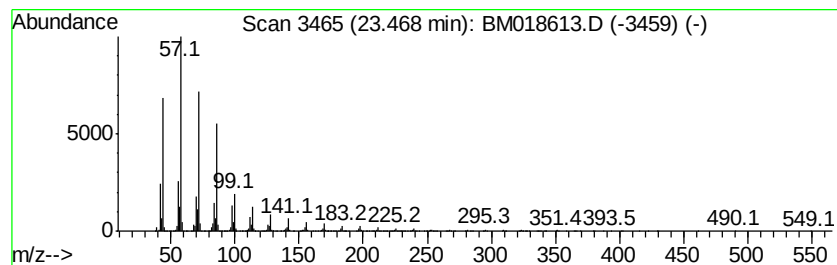
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tricosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.47	33.04 ng	6150500	Perylene-d12	23.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	94
2		Tetracosane	338	C24H50	000646-31-1	94
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
4		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	91
5		Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011719\
Data File : BM018613.D
Acq On : 17 Jan 2019 14:36
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Sample : K1147-22
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ClientSampleId :
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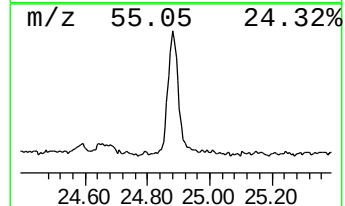
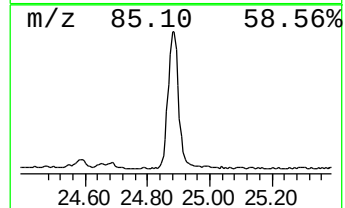
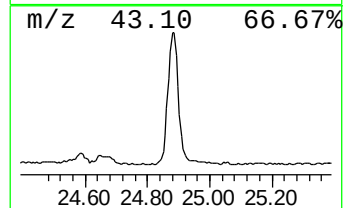
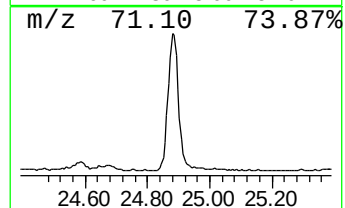
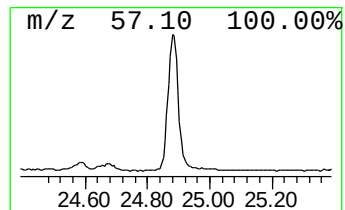
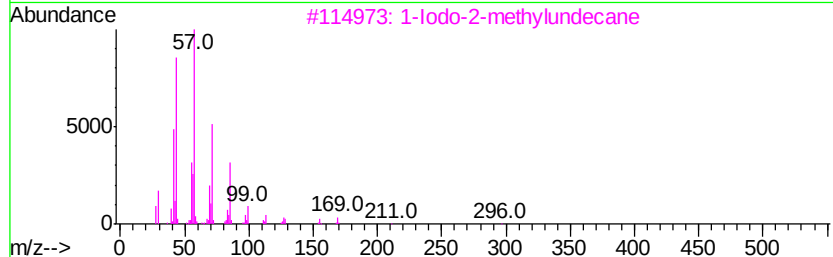
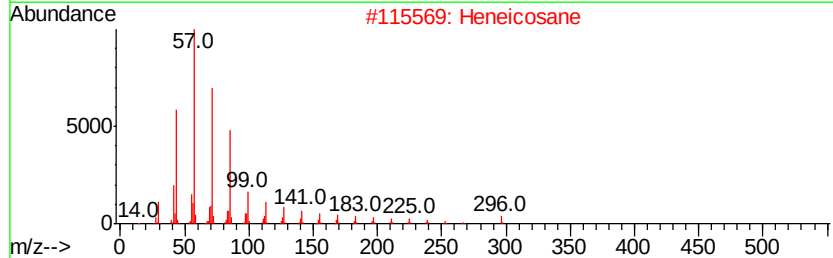
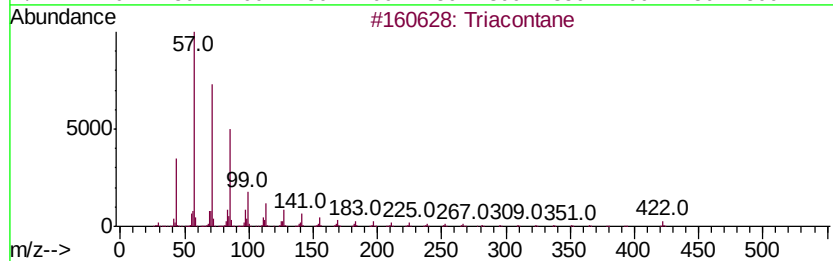
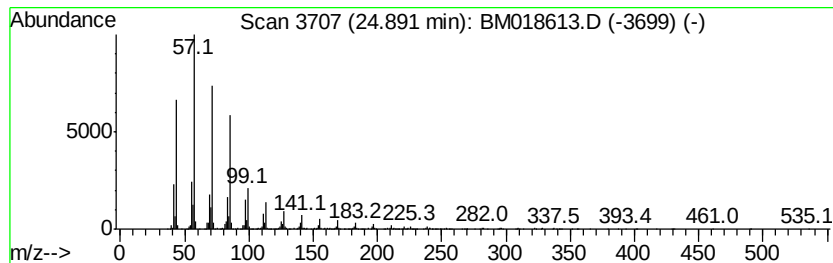
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Triacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.89	12.79 ng	2380690	Perylene-d12	23.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM011719\
Data File : BM018613.D
Acq On : 17 Jan 2019 14:36
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Sample : K1147-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM010919.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
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2-Pentanone, 4-hy...	4.89	4.6	ng	504902	1	7.74	2176120	20.0
unknown7.27	7.27	110.9	ng	12070400	1	7.74	2176120	20.0
n-Hexadecanoic acid	18.03	11.3	ng	2042460	4	17.14	3609190	20.0
Octadecanoic acid	19.31	6.4	ng	1227620	5	21.33	3844640	20.0
Ethanol, 2-(hexad...	21.03	3.5	ng	675941	5	21.33	3844640	20.0
Nonacosane	21.47	4.5	ng	857920	5	21.33	3844640	20.0
Octadecane	21.91	12.4	ng	2390350	5	21.33	3844640	20.0
Heptadecane, 9-he...	22.37	24.9	ng	4776090	5	21.33	3844640	20.0
Tetracosane	22.89	32.6	ng	6079170	6	23.60	3723430	20.0
Tricosane	23.47	33.0	ng	6150500	6	23.60	3723430	20.0
Triacontane	24.89	12.8	ng	2380690	6	23.60	3723430	20.0