

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043899.D
 Acq On : 3 Jan 2020 22:32
 Operator : JU
 Sample : K6449-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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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_G\METHODS\8270-BG123019.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	3.146	5	10	34	rVB	1892551	2873365	30.45%	5.150%
2	5.067	331	337	345	rBV	282652	425801	4.51%	0.763%
3	5.538	409	417	426	rBV	1390145	2175544	23.05%	3.899%
4	7.124	680	687	698	rVB	980963	1732886	18.36%	3.106%
5	7.494	741	750	760	rBV	2200555	3818218	40.46%	6.843%
6	7.964	823	830	838	rBV	437212	766563	8.12%	1.374%
7	8.287	876	885	893	rBV	1657459	2866862	30.38%	5.138%
8	9.116	1017	1026	1037	rBV	1431902	2544939	26.97%	4.561%
9	10.761	1298	1306	1314	rBV	621987	1151496	12.20%	2.064%
10	13.217	1717	1724	1733	rBV	3643259	5995289	63.53%	10.745%
11	14.045	1857	1865	1878	rBV	268717	403992	4.28%	0.724%
12	14.592	1949	1958	1966	rBV	1099148	1730556	18.34%	3.101%
13	16.078	2204	2211	2229	rBV	3380612	5139568	54.46%	9.211%
14	17.335	2418	2425	2436	rBV	1409868	2147717	22.76%	3.849%
15	17.459	2438	2446	2454	rVB2	133941	210709	2.23%	0.378%
16	19.950	2864	2870	2883	rBV	6756694	9436724	100.00%	16.912%
17	20.273	2921	2925	2929	rVB	168715	191881	2.03%	0.344%
18	20.767	3002	3009	3013	rBV	246491	299576	3.17%	0.537%
19	21.254	3087	3092	3106	rBV	1233005	1943644	20.60%	3.483%
20	21.583	3142	3148	3156	rVB	1365612	2206752	23.38%	3.955%
21	21.807	3181	3186	3192	rBV	288152	412277	4.37%	0.739%
22	22.406	3282	3288	3301	rBV2	395652	726731	7.70%	1.302%
23	23.088	3397	3404	3415	rBV	454404	869010	9.21%	1.557%
24	23.875	3530	3538	3550	rBV	440879	1007730	10.68%	1.806%
25	24.709	3670	3680	3689	rBV2	823904	2277639	24.14%	4.082%
26	24.803	3689	3696	3713	rVB2	380551	986741	10.46%	1.768%
27	25.908	3875	3884	3894	rBV2	264860	803066	8.51%	1.439%
28	26.014	3896	3902	3914	rVB5	86128	241655	2.56%	0.433%
29	27.242	4101	4111	4120	rBV3	123846	411622	4.36%	0.738%

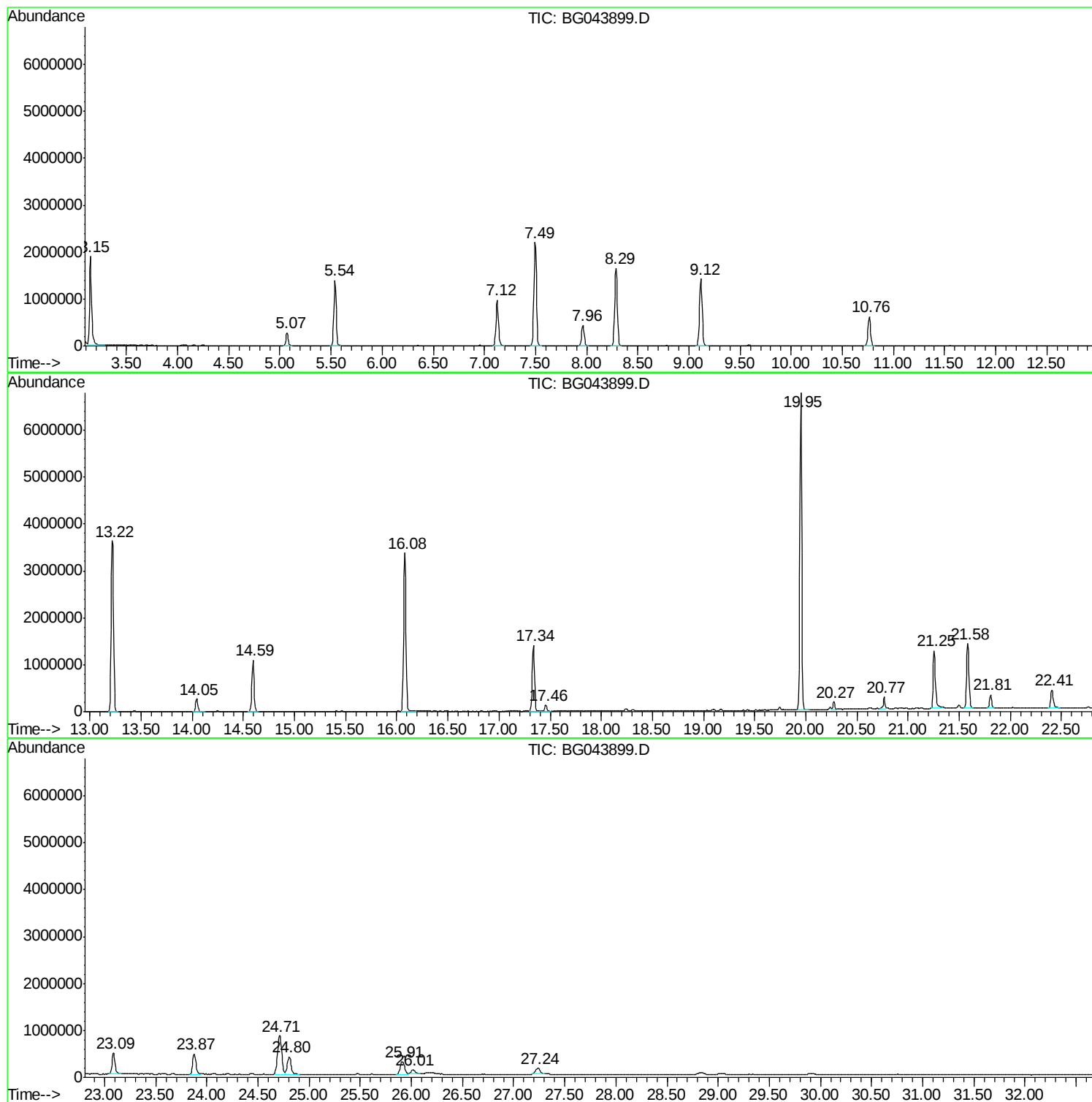
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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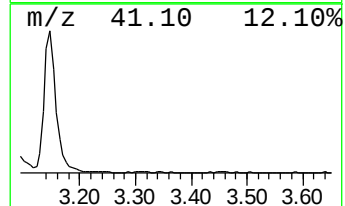
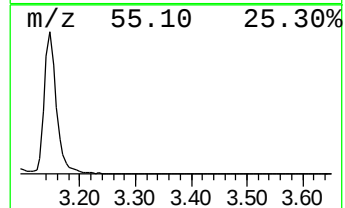
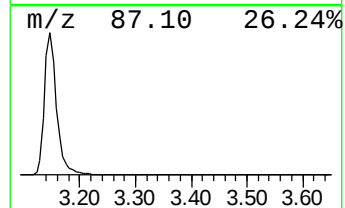
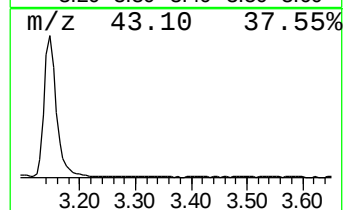
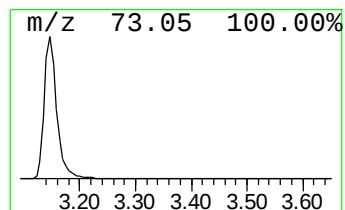
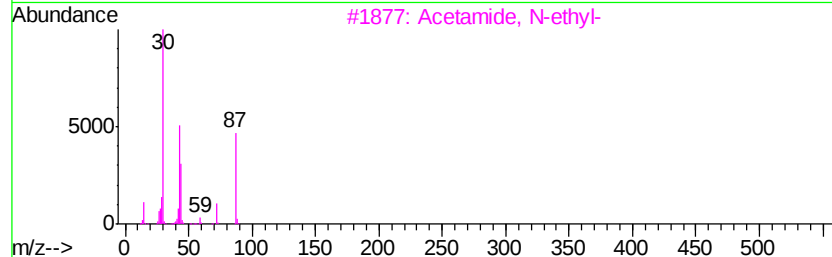
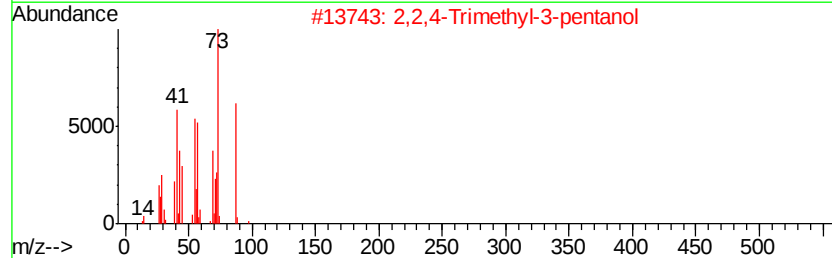
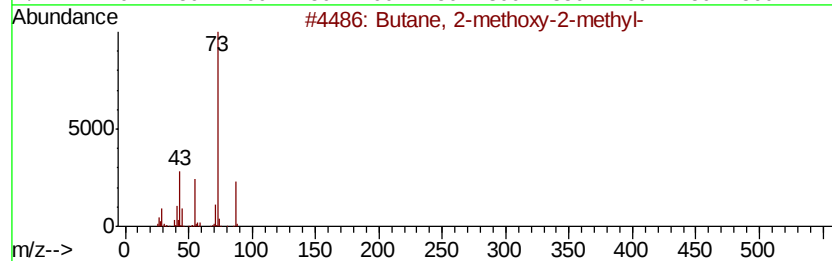
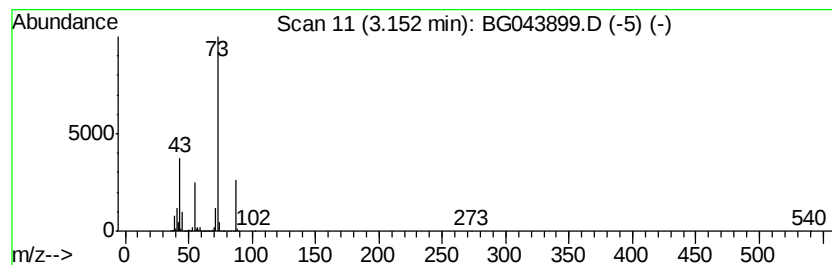
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	74.97 ng	2873370	1,4-Dichlorobenzene-d4	7.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	35
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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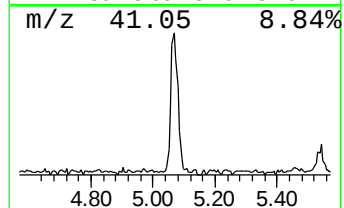
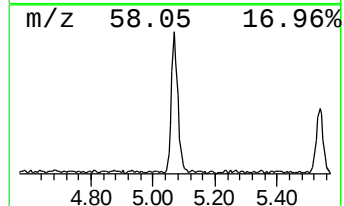
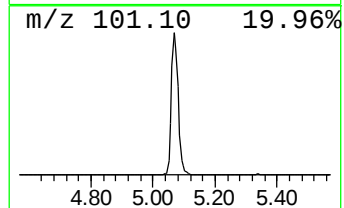
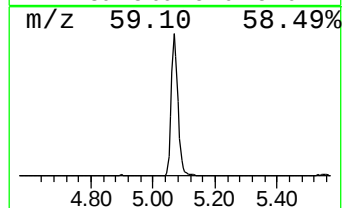
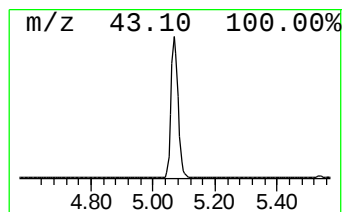
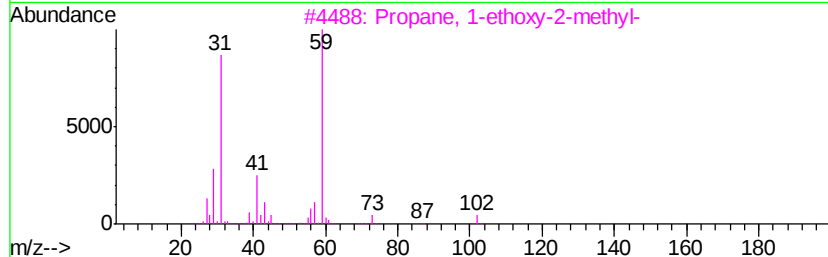
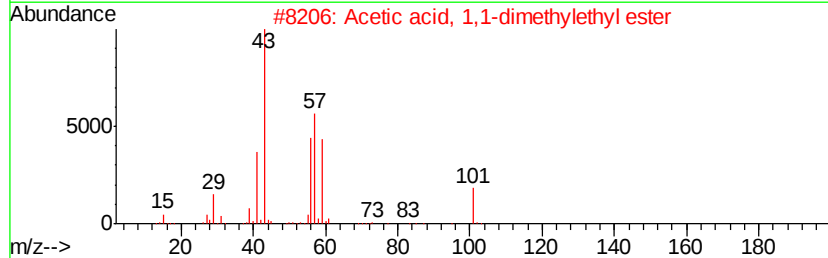
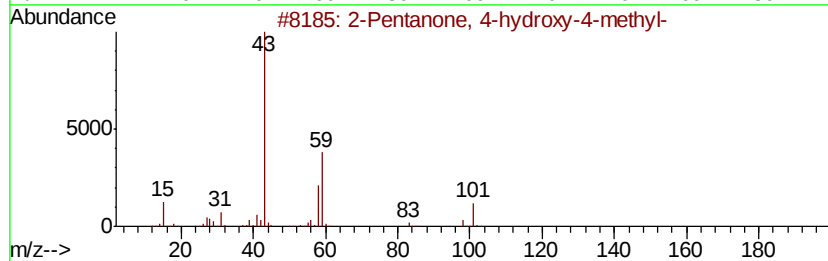
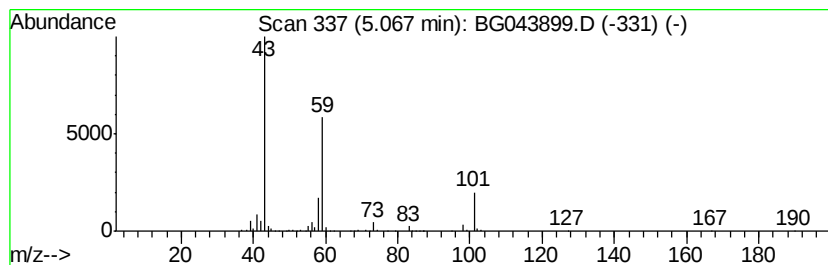
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	11.11 ng	425801	1,4-Dichlorobenzene-d4	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	32
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	23
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17



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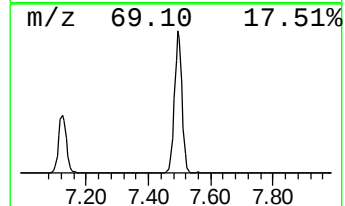
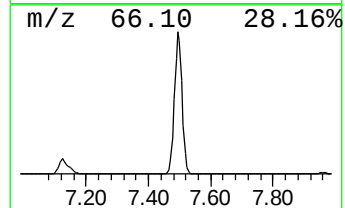
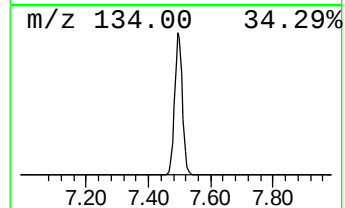
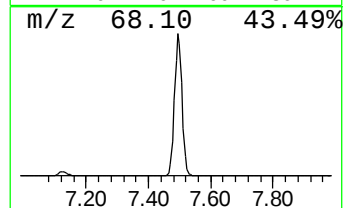
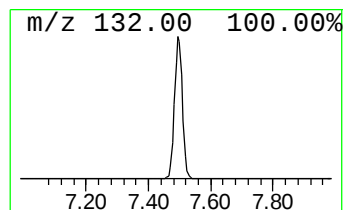
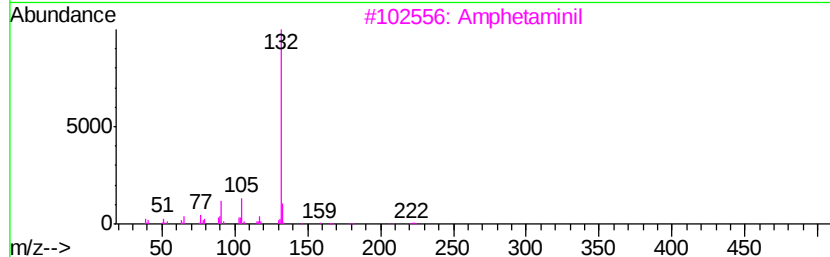
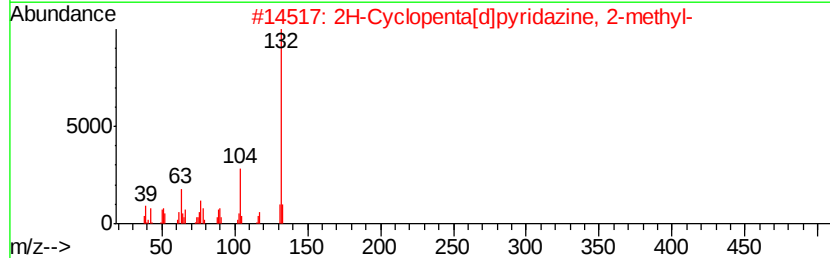
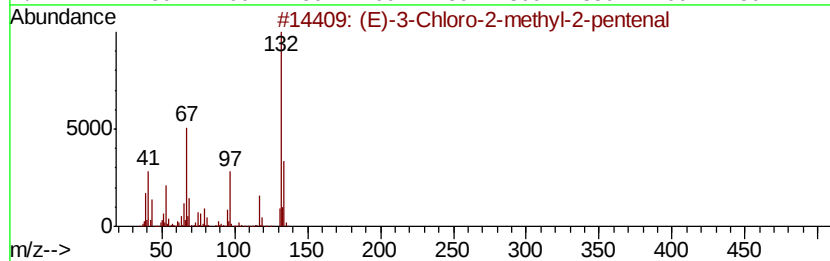
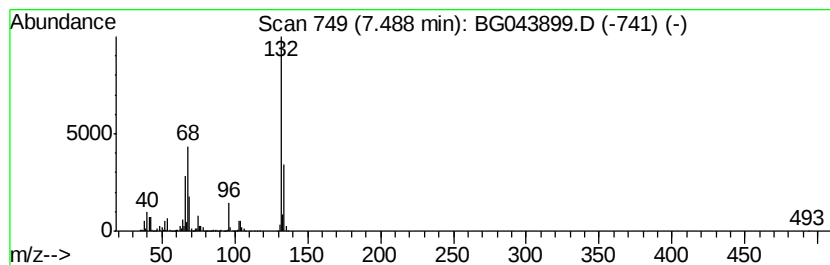
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Peak Number 3 unknown7.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.49	99.62 ng	3818220	1,4-Dichlorobenzene-d4	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	12
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		1-Isoquinolinemethanol, .alpha.-...	191	C12H17NO	114608-92-3	12
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	12



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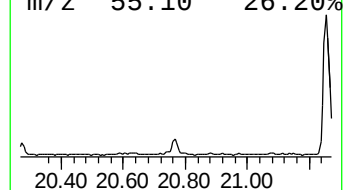
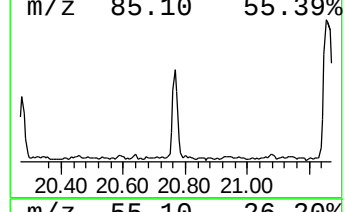
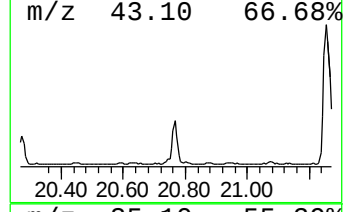
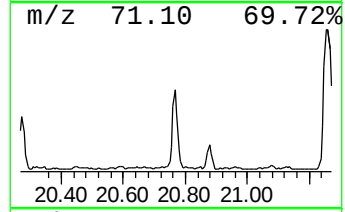
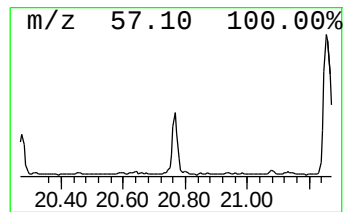
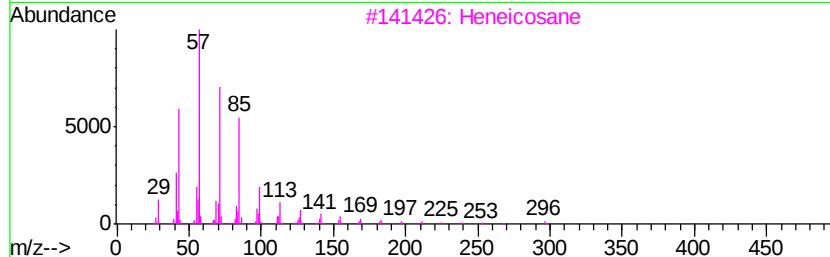
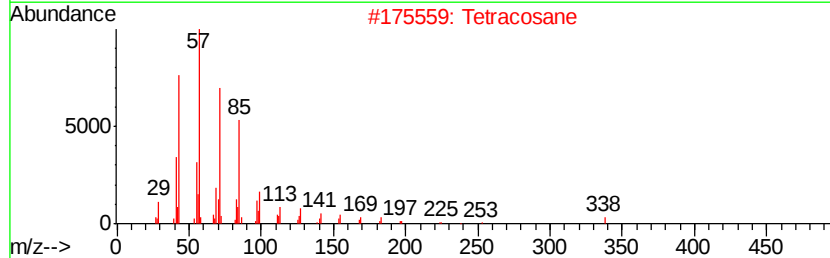
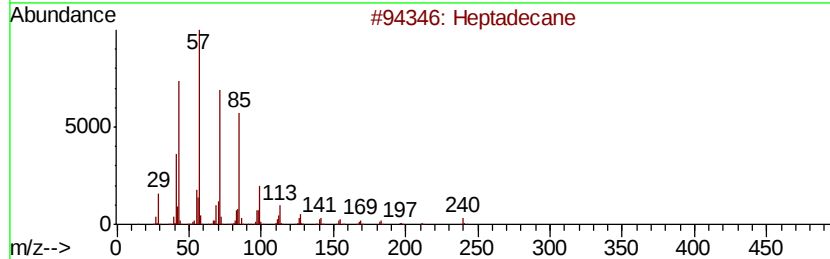
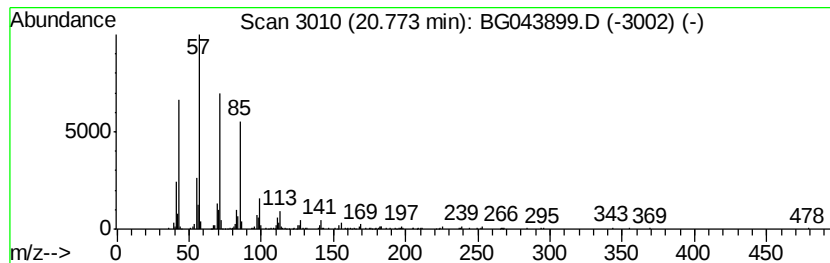
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Peak Number 5 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.77	2.72 ng	299576	Chrysene-d12	21.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Tetracosane		338	C24H50	000646-31-1	91
3	Heneicosane		296	C21H44	000629-94-7	91
4	Eicosane		282	C20H42	000112-95-8	91
5	Hexadecane		226	C16H34	000544-76-3	87



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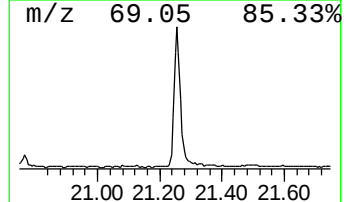
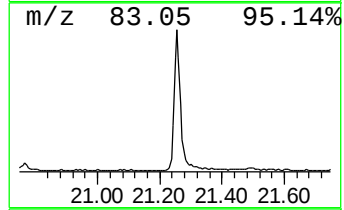
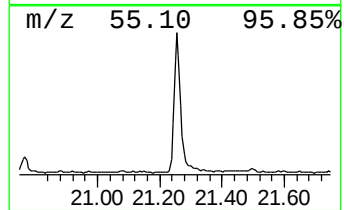
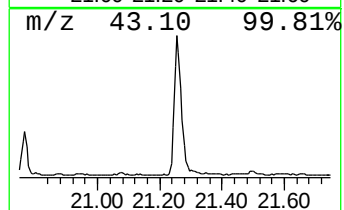
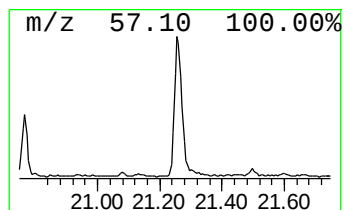
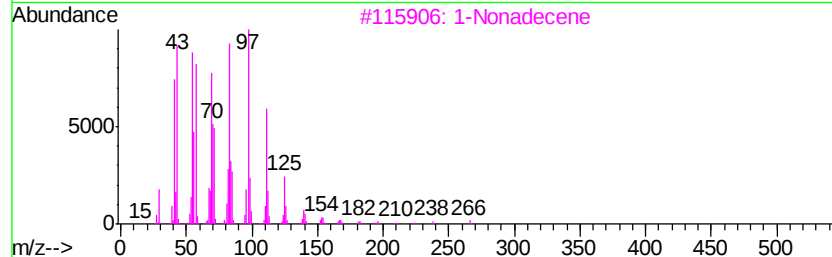
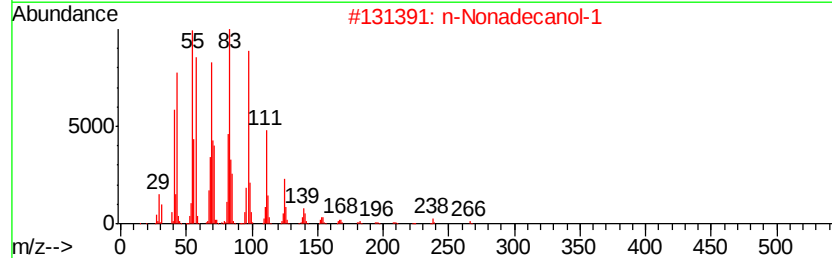
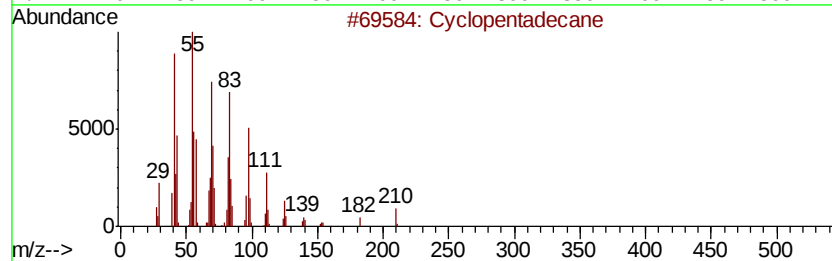
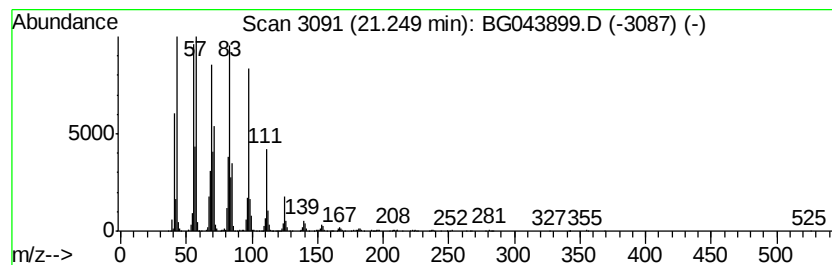
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 Peak Number 6 Cyclopentadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.25	17.62 ng	1943640	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	93
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	91
3		1-Nonadecene	266	C19H38	018435-45-5	91
4		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
5		Behenic alcohol	326	C22H46O	000661-19-8	91



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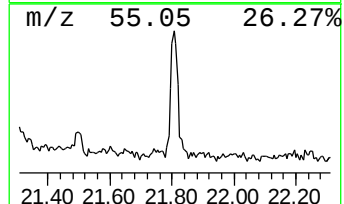
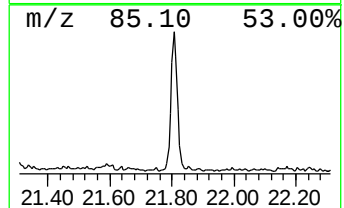
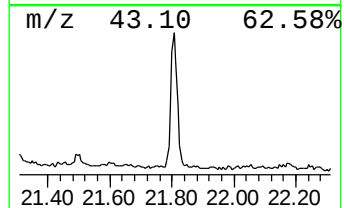
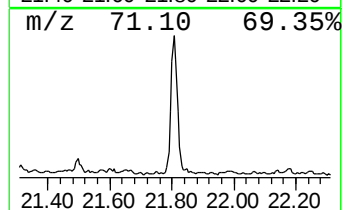
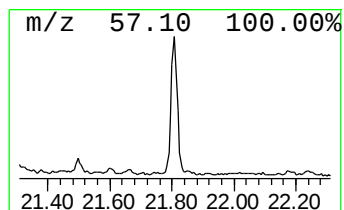
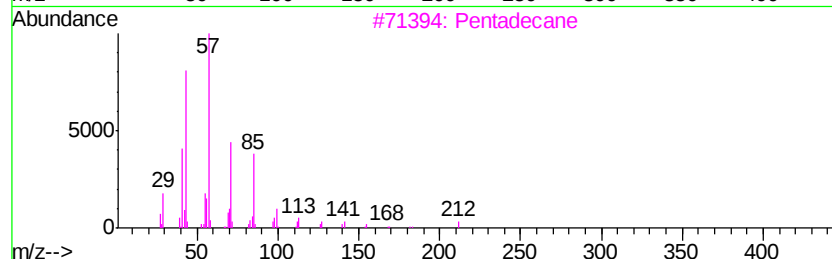
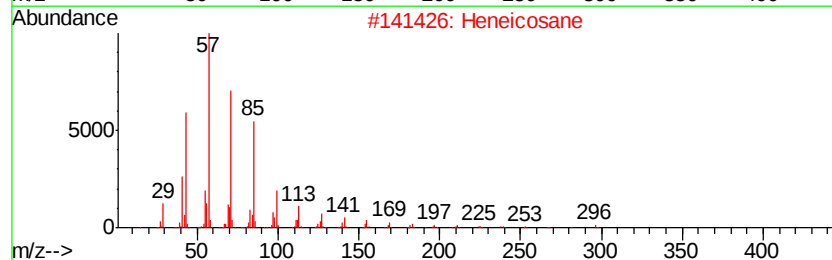
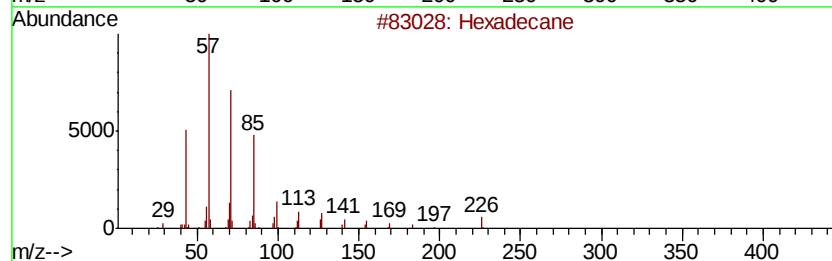
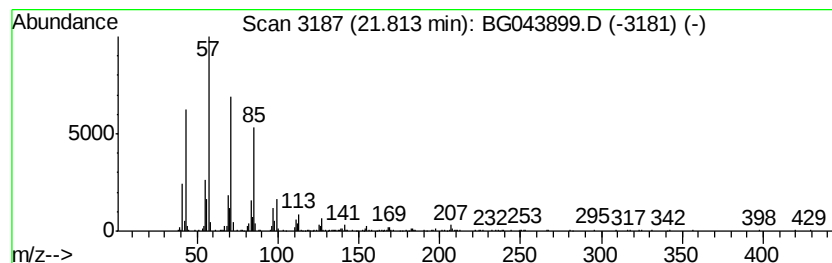
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	3.74 ng	412277	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Pentadecane	212	C15H32	000629-62-9	90
4		Eicosane	282	C20H42	000112-95-8	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043899.D
Acq On : 3 Jan 2020 22:32
Operator : JU
Sample : K6449-10
Misc :
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Instrument :
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ClientSampleId :
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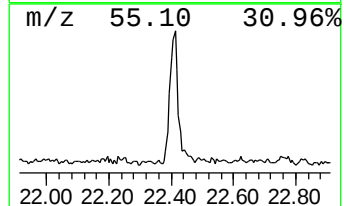
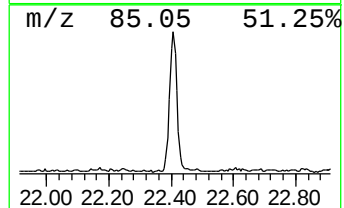
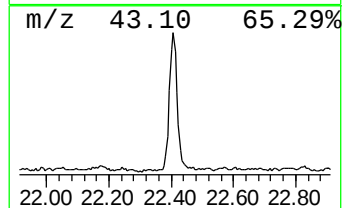
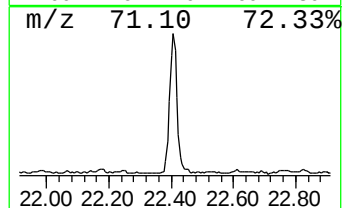
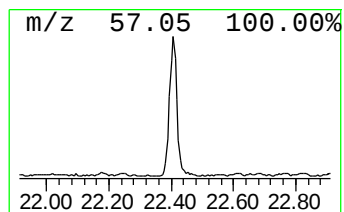
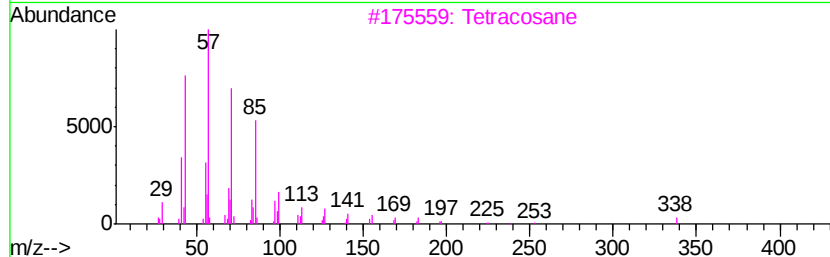
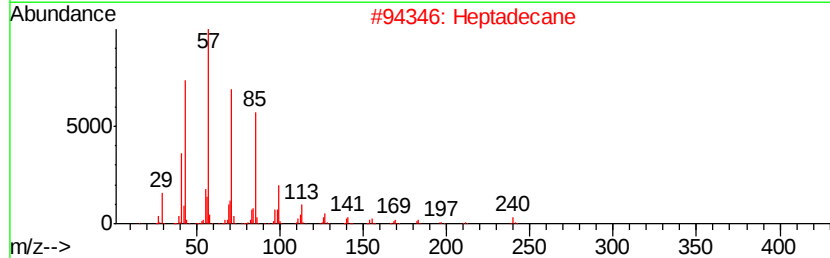
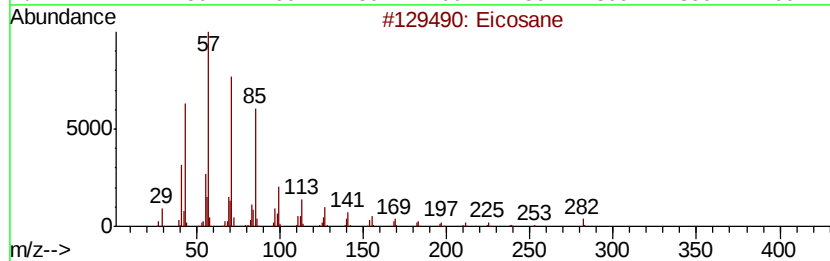
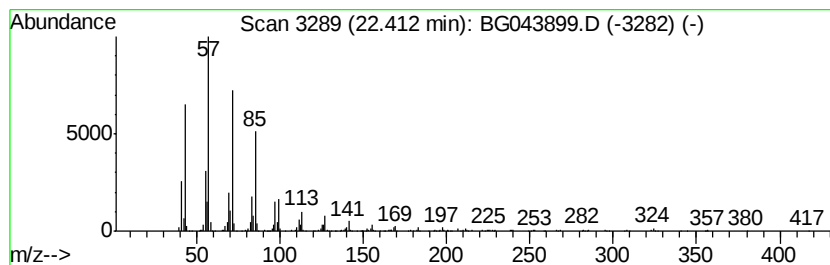
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.41	6.59 ng	726731	Chrysene-d12	21.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Heptadecane		240	C17H36	000629-78-7	95
3	Tetracosane		338	C24H50	000646-31-1	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Octacosane		394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
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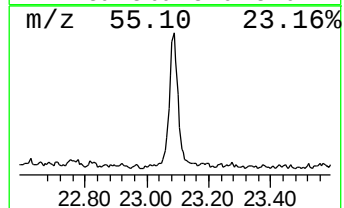
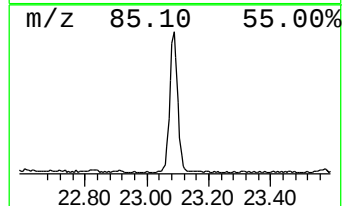
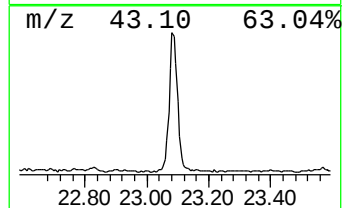
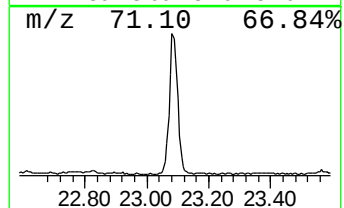
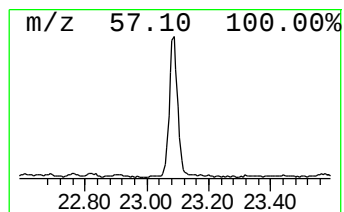
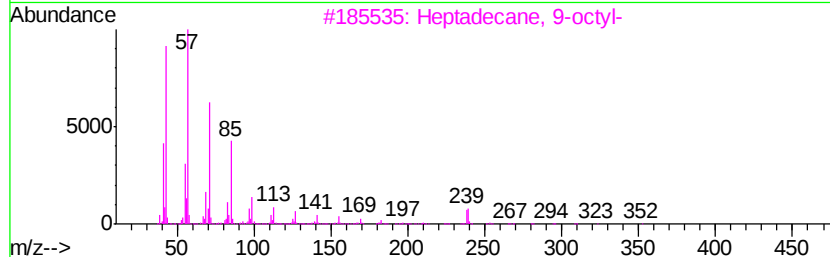
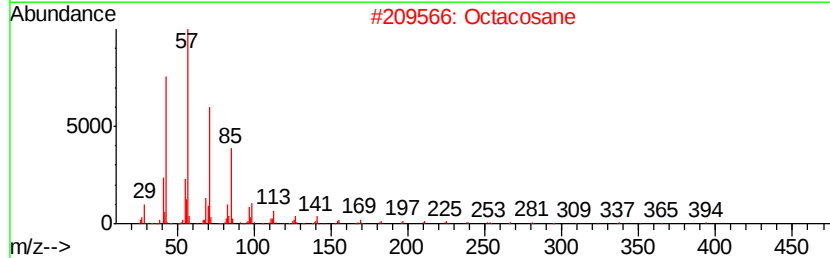
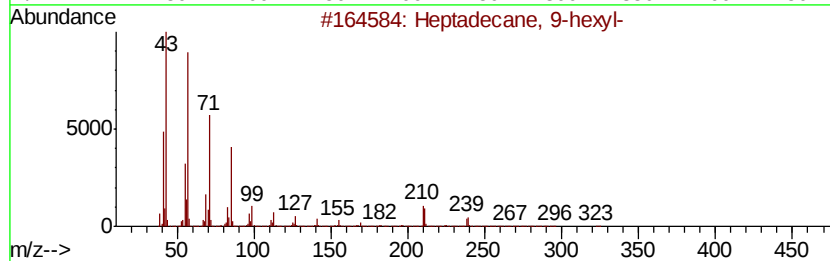
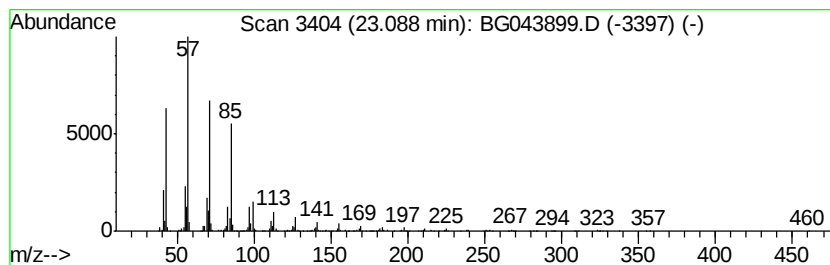
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane, 9-hexyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.09	7.88 ng	869010	Chrysene-d12	21.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-hexyl-		324	C23H48	055124-79-3	94
2	Octacosane		394	C28H58	000630-02-4	91
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Tetracosane		338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043899.D
Acq On : 3 Jan 2020 22:32
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Misc :
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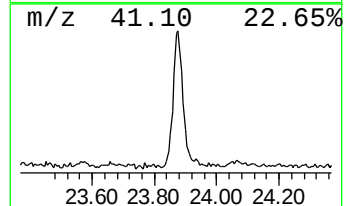
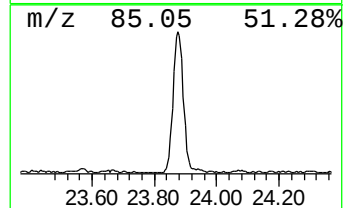
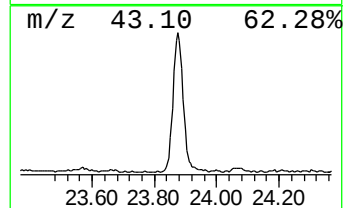
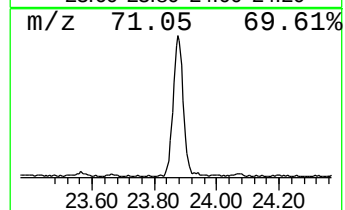
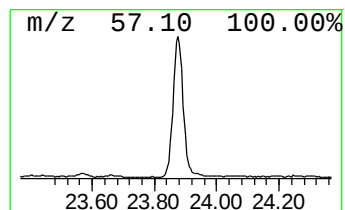
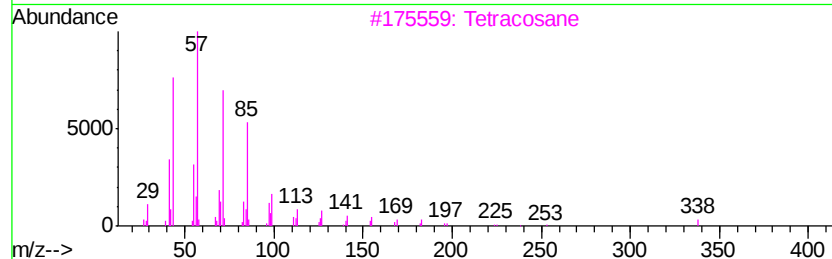
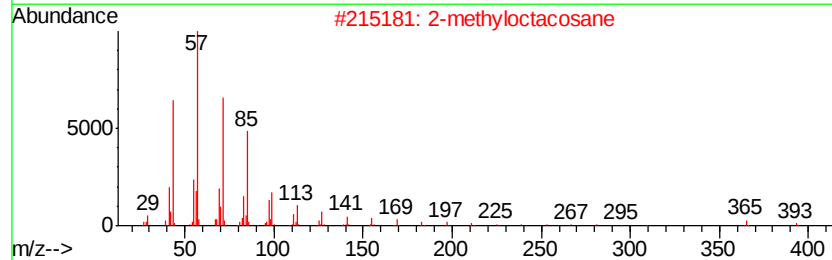
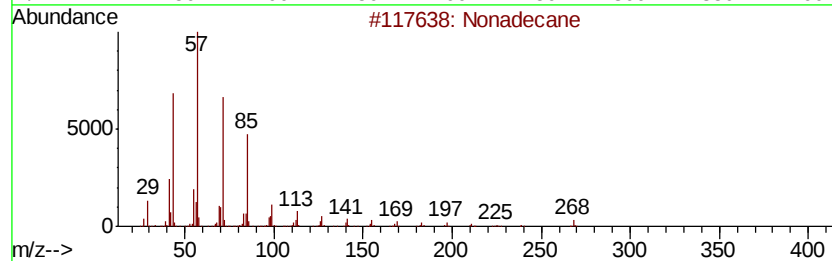
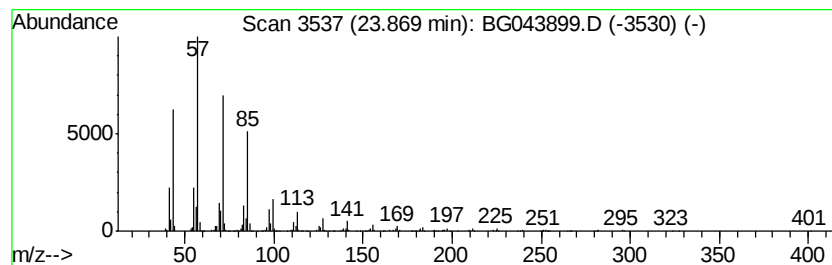
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.87	8.85 ng	1007730	Perylene-d12	24.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			2-methyloctacosane	408	C29H60	1000376-72-8	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043899.D
Acq On : 3 Jan 2020 22:32
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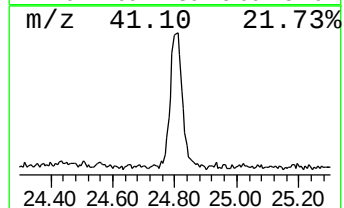
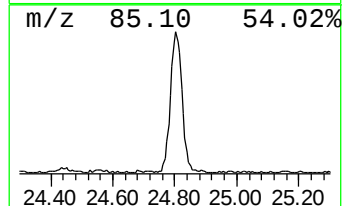
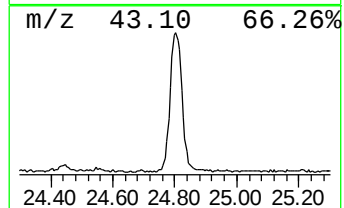
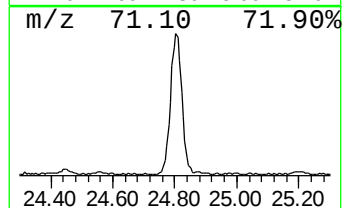
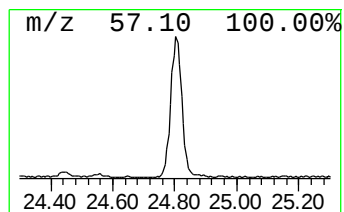
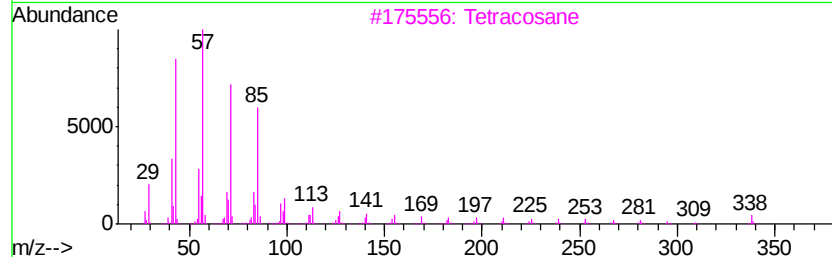
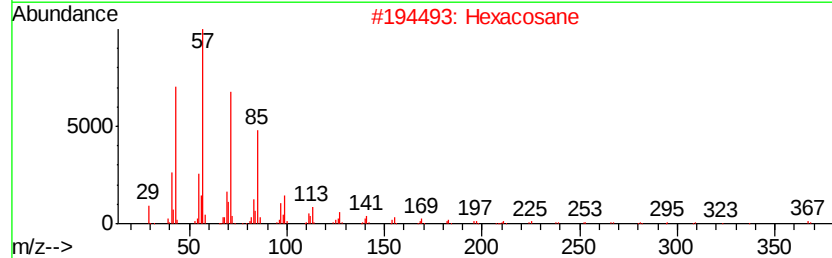
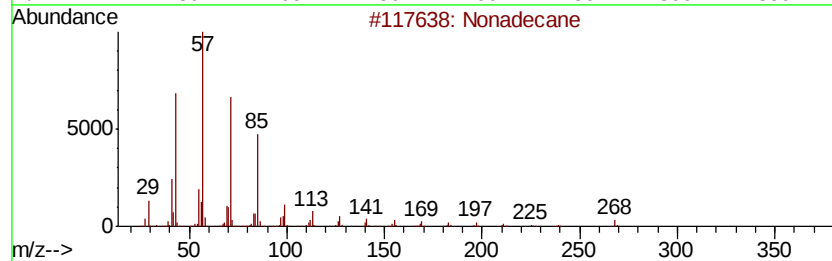
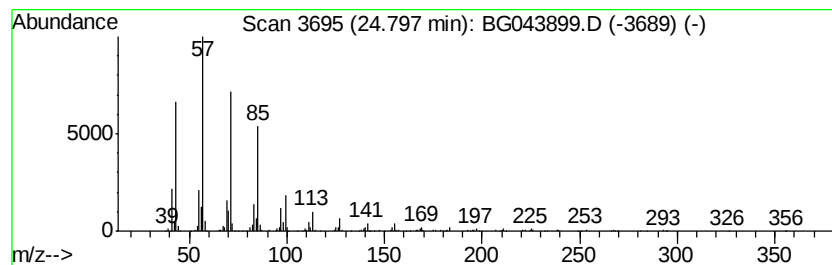
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.80	8.66 ng	986741	Perylene-d12	24.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Hexacosane	366	C26H54	000630-01-3	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heneicosane	296	C21H44	000629-94-7	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043899.D
 Acq On : 3 Jan 2020 22:32
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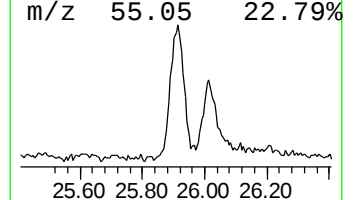
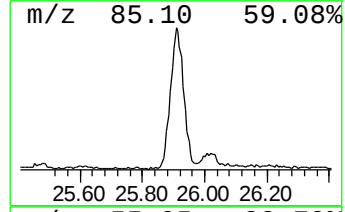
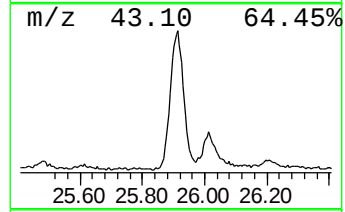
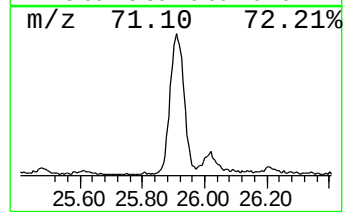
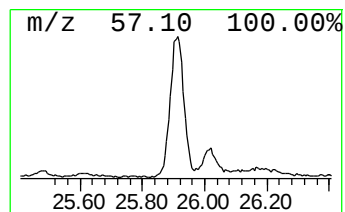
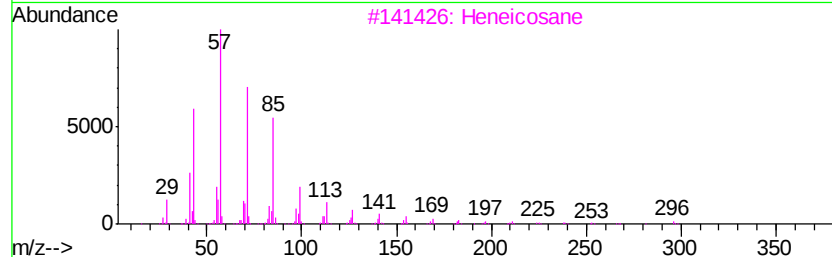
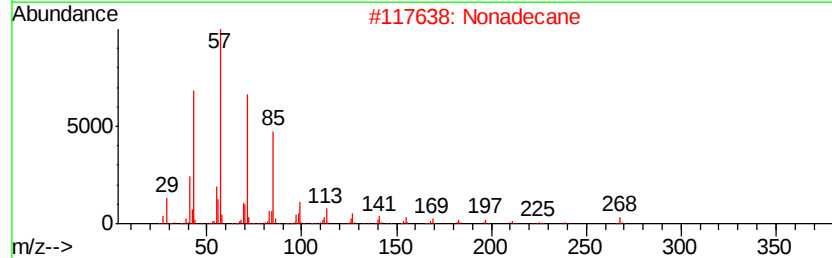
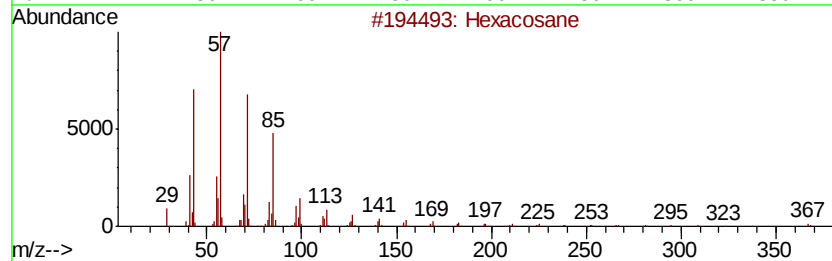
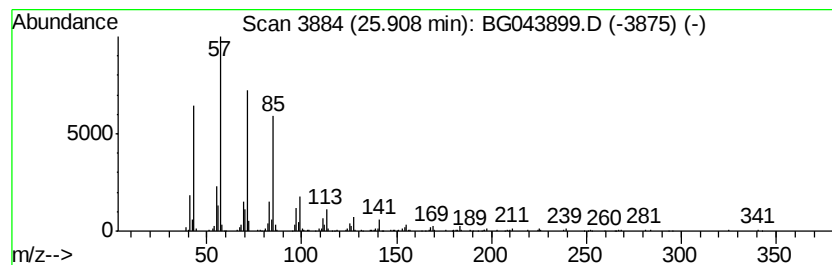
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Heptadecane, 9-octyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.91	7.05 ng	803066	Perylene-d12	24.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	94
2			Nonadecane	268	C19H40	000629-92-5	93
3			Heneicosane	296	C21H44	000629-94-7	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043899.D
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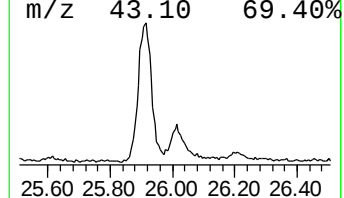
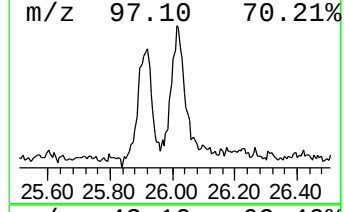
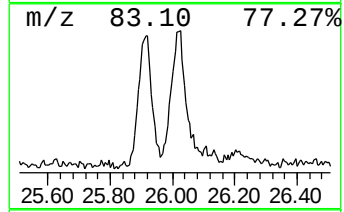
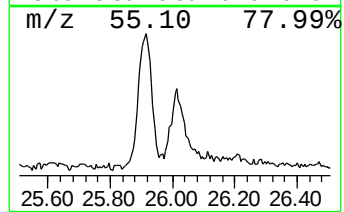
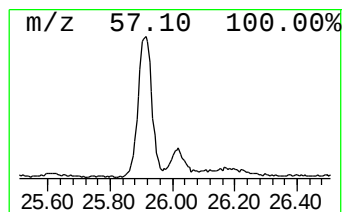
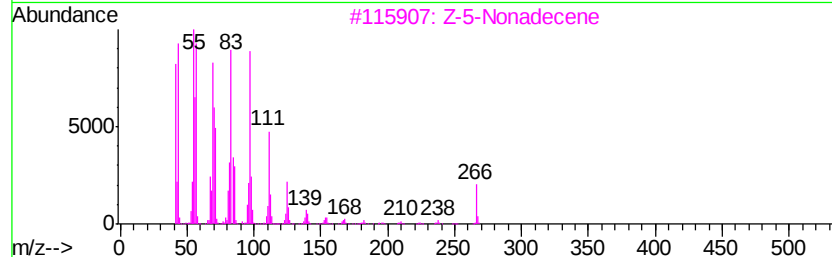
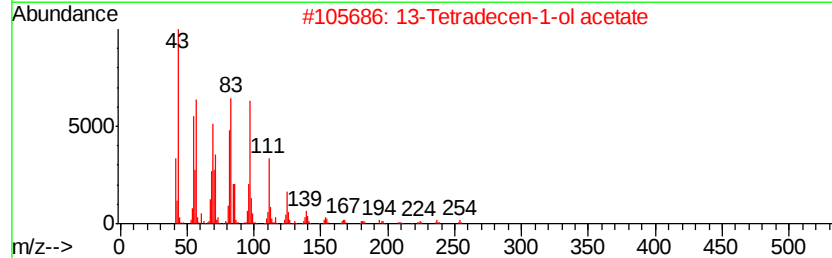
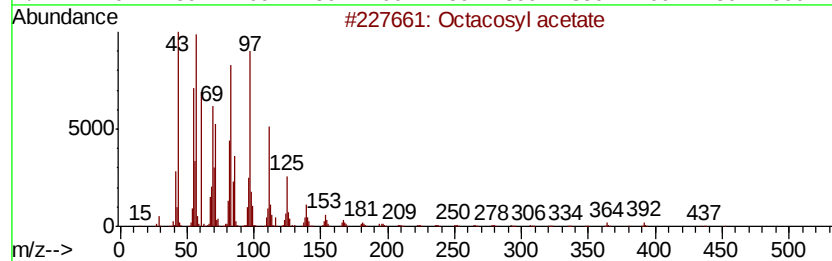
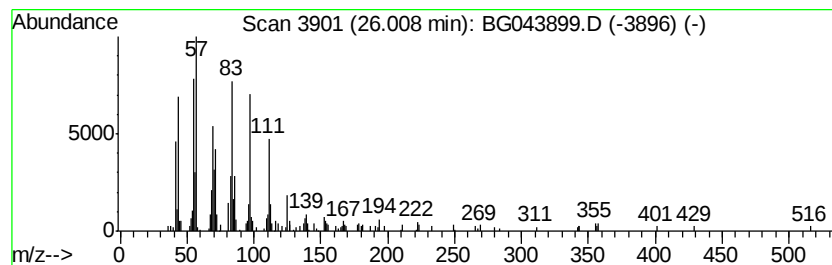
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Octacosyl acetate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.01	2.12 ng	241655	Perylene-d12	24.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosyl acetate	452	C30H60O2	018206-97-8	90
2		13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	87
3		Z-5-Nonadecene	266	C19H38	1000131-11-8	83
4		1-Docosene	308	C22H44	001599-67-3	81
5		Octatriacontyl trifluoroacetate	647	C40H77F3O2	1000351-88-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
Data File : BG043899.D
Acq On : 3 Jan 2020 22:32
Operator : JU
Sample : K6449-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

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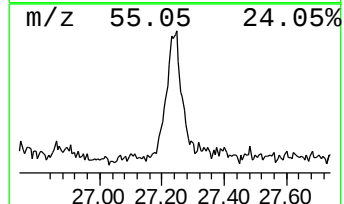
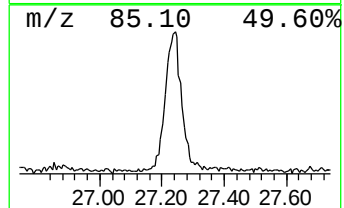
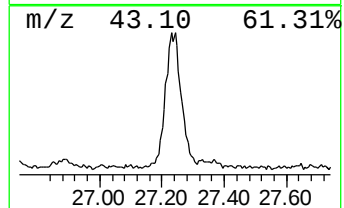
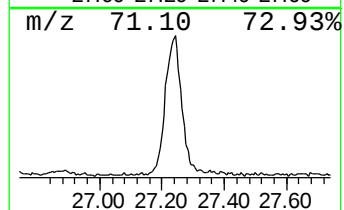
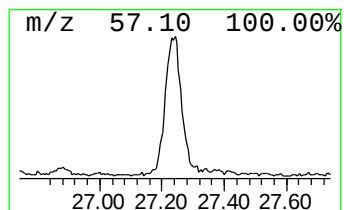
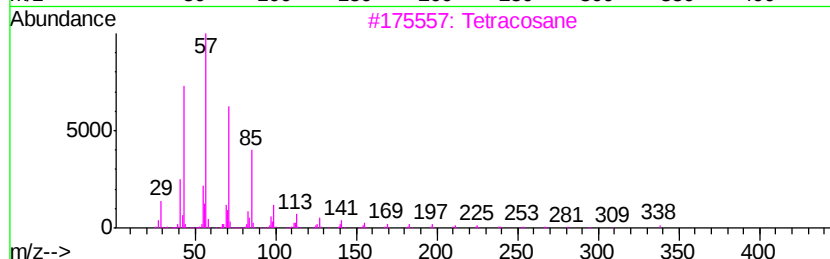
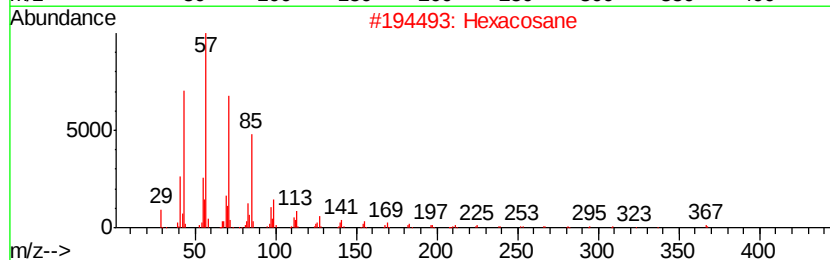
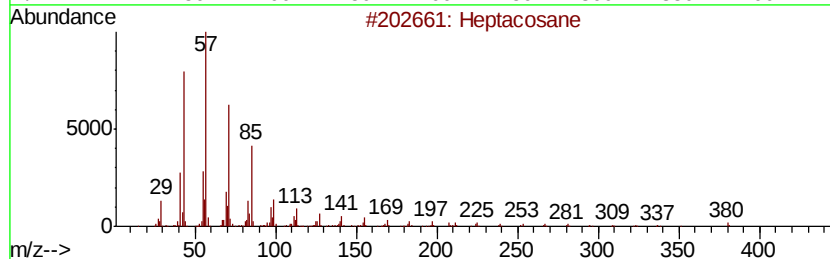
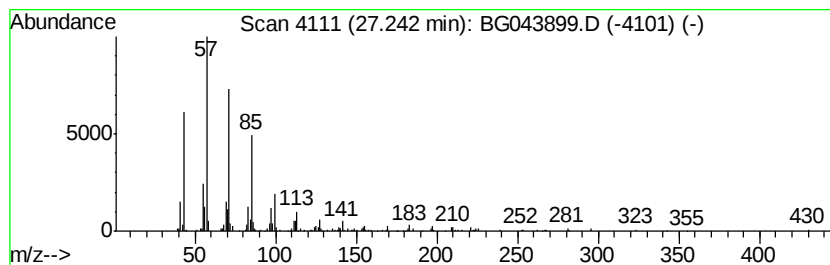
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.24	3.61 ng	411622	Perylene-d12	24.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	91
2			Hexacosane	366	C26H54	000630-01-3	90
3			Tetracosane	338	C24H50	000646-31-1	87
4			Octacosane	394	C28H58	000630-02-4	87
5			Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG010320\
 Data File : BG043899.D
 Acq On : 3 Jan 2020 22:32
 Operator : JU
 Sample : K6449-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FIELD-BLANK

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.15	75.0	ng	2873370	1	7.96	766563	20.0
2-Pentanone, 4-hy...	5.07	11.1	ng	425801	1	7.96	766563	20.0
unknown7.49	7.49	99.6	ng	3818220	1	7.96	766563	20.0
Heptadecane	20.77	2.7	ng	299576	5	21.58	2206750	20.0
Cyclopentadecane	21.25	17.6	ng	1943640	5	21.58	2206750	20.0
Hexadecane	21.81	3.7	ng	412277	5	21.58	2206750	20.0
Eicosane	22.41	6.6	ng	726731	5	21.58	2206750	20.0
Heptadecane, 9-he...	23.09	7.9	ng	869010	5	21.58	2206750	20.0
Nonadecane	23.87	8.8	ng	1007730	6	24.71	2277640	20.0
Hexacosane	24.80	8.7	ng	986741	6	24.71	2277640	20.0
Heptadecane, 9-oc...	25.91	7.0	ng	803066	6	24.71	2277640	20.0
Octacosyl acetate	26.01	2.1	ng	241655	6	24.71	2277640	20.0
Heptacosane	27.24	3.6	ng	411622	6	24.71	2277640	20.0