

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
 Data File : BG044024.D
 Acq On : 13 Jan 2020 22:57
 Operator : JU
 Sample : L1074-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CIY1-GW-102-13.5

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	3.124	3	6	28	rVB	1478966	2481478	33.92%	5.555%
2	5.051	327	334	343	rBV	178011	278477	3.81%	0.623%
3	5.521	407	414	428	rBV	819706	1347407	18.42%	3.016%
4	7.108	677	684	703	rBV	523312	994730	13.60%	2.227%
5	7.478	739	747	764	rBV	1591311	2811616	38.44%	6.294%
6	7.948	819	827	834	rBV	369091	637383	8.71%	1.427%
7	8.271	873	882	890	rBV	1419020	2523112	34.49%	5.648%
8	9.100	1013	1023	1034	rBV	1158223	2097152	28.67%	4.695%
9	10.745	1294	1303	1318	rBV	479166	904335	12.36%	2.024%
10	13.201	1713	1721	1732	rBV	2863587	4705736	64.33%	10.534%
11	14.029	1856	1862	1870	rBV	220249	344052	4.70%	0.770%
12	14.576	1947	1955	1969	rBV	835837	1278042	17.47%	2.861%
13	16.062	2201	2208	2221	rBV	2562653	3963428	54.18%	8.873%
14	17.319	2415	2422	2432	rBV	1025884	1539325	21.04%	3.446%
15	19.934	2861	2867	2879	rBV	4766581	7314974	100.00%	16.375%
16	21.238	3083	3089	3102	rBV	785482	1246093	17.03%	2.790%
17	21.567	3139	3145	3156	rBV	955668	1647292	22.52%	3.688%
18	21.785	3177	3182	3188	rVB	178605	245875	3.36%	0.550%
19	22.378	3278	3283	3298	rVB2	289535	566681	7.75%	1.269%
20	23.060	3393	3399	3407	rVB	362965	661071	9.04%	1.480%
21	23.236	3423	3429	3433	rBV	167309	337431	4.61%	0.755%
22	23.847	3526	3533	3542	rBV	369587	786334	10.75%	1.760%
23	24.323	3605	3614	3624	rBV	163822	429154	5.87%	0.961%
24	24.687	3666	3676	3685	rBV2	469959	1624792	22.21%	3.637%
25	24.775	3685	3691	3703	rVB2	333427	896461	12.26%	2.007%
26	25.621	3824	3835	3853	rVB3	191320	705846	9.65%	1.580%
27	25.880	3870	3879	3890	rBV	218703	673114	9.20%	1.507%
28	27.225	4095	4108	4133	rVB9	257617	1347881	18.43%	3.017%
29	29.276	4446	4457	4458	rBV2	115005	281025	3.84%	0.629%

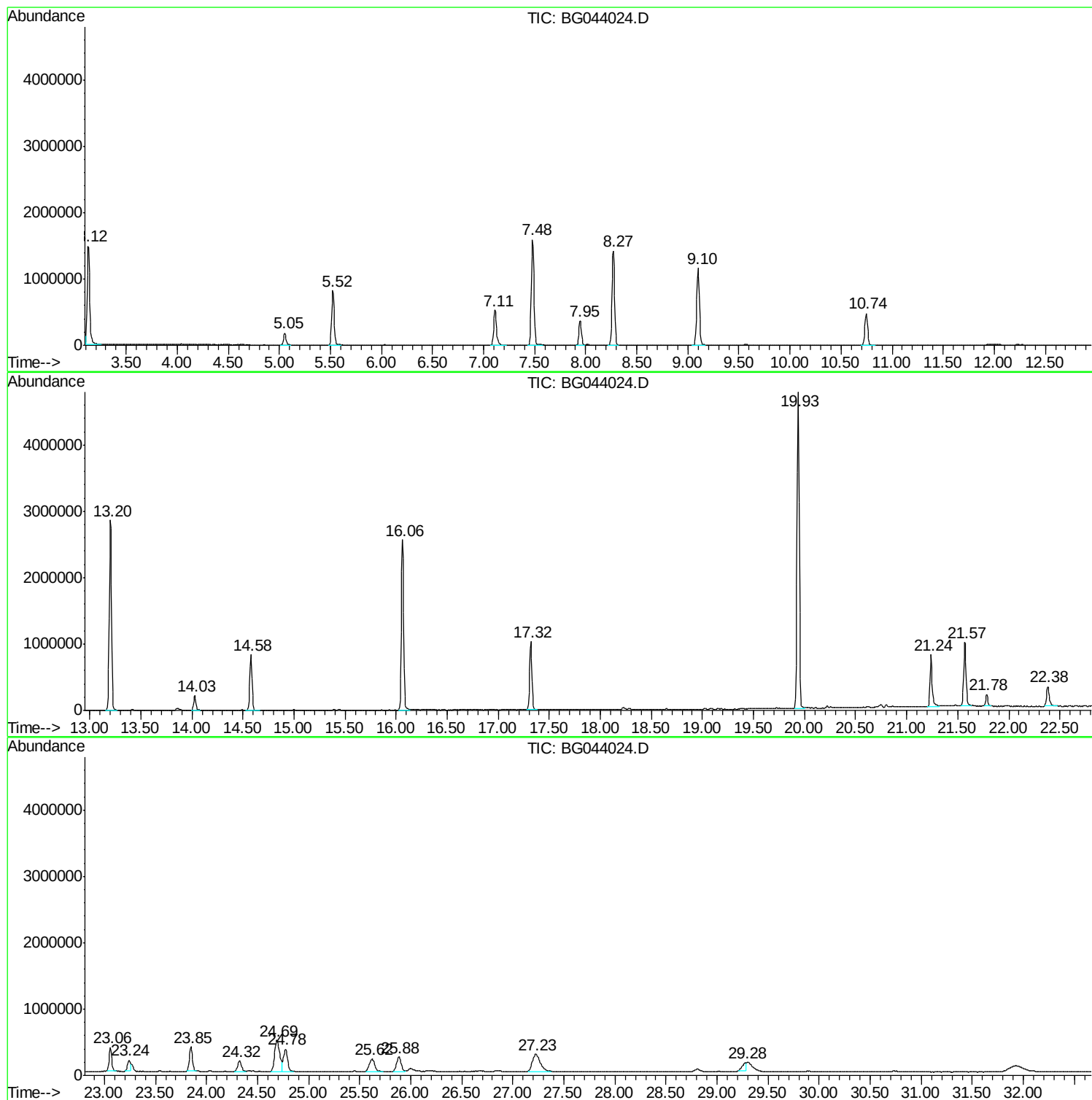
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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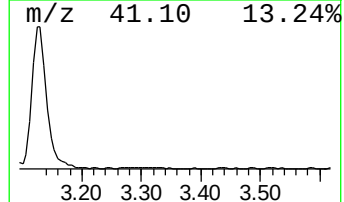
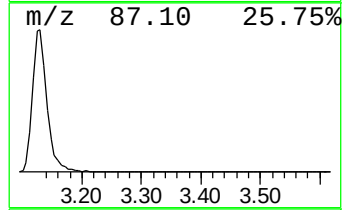
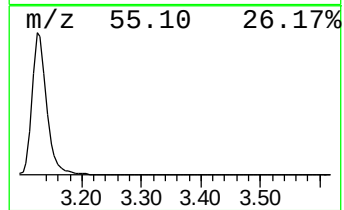
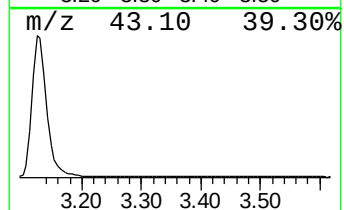
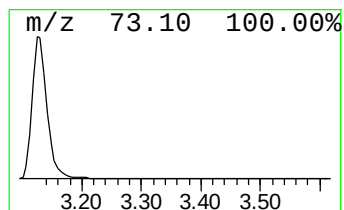
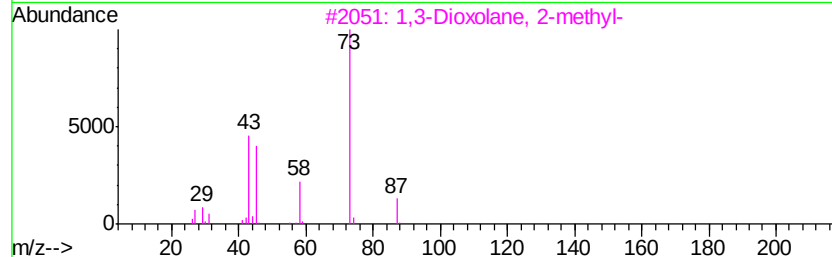
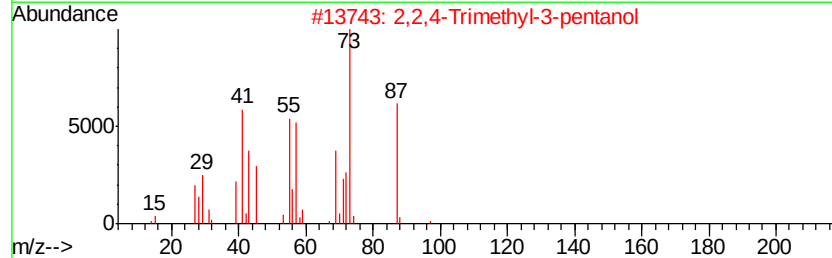
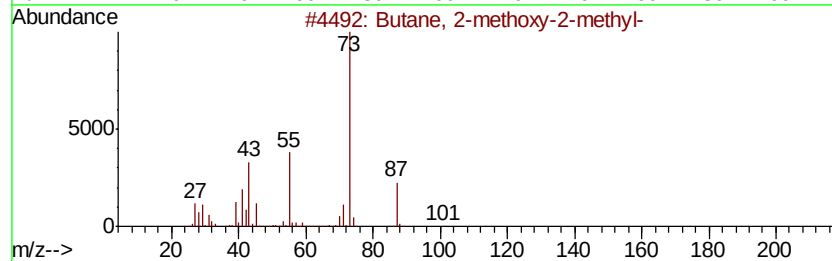
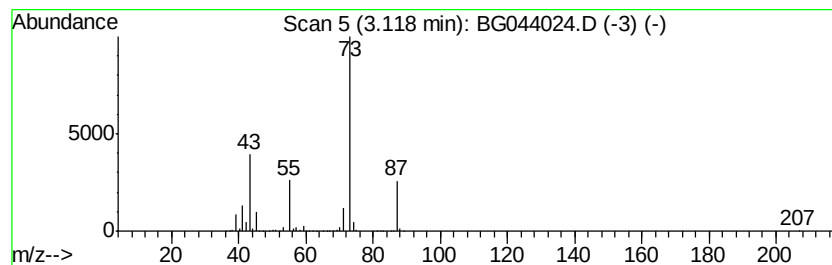
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.12	77.86 ng	2481480	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22



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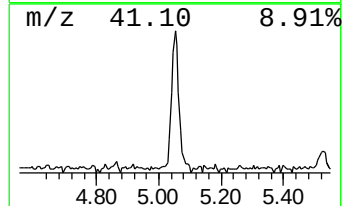
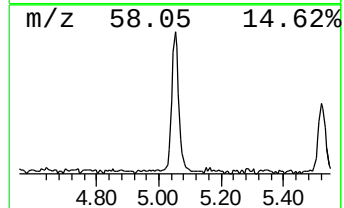
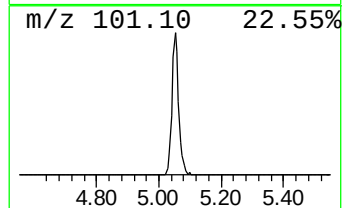
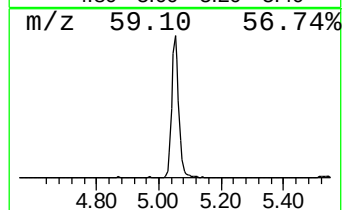
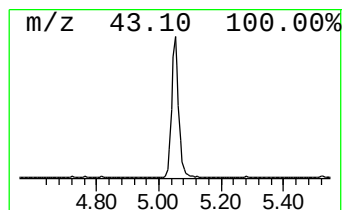
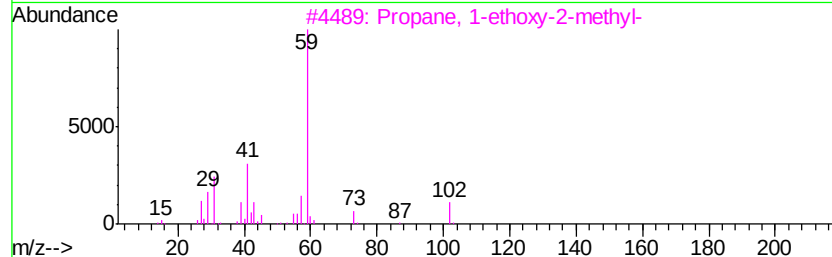
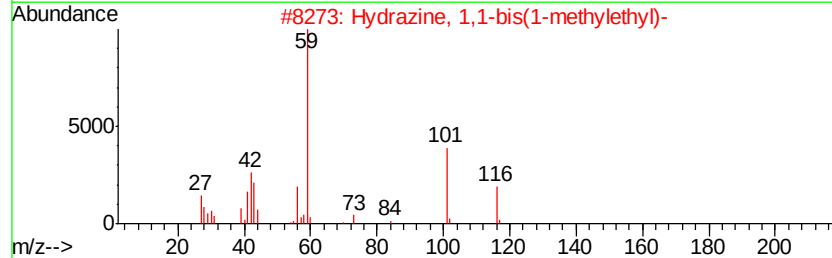
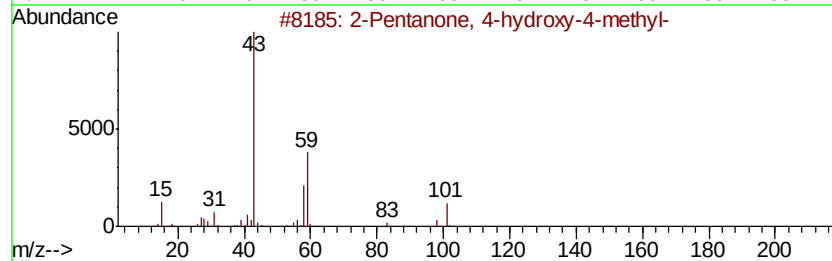
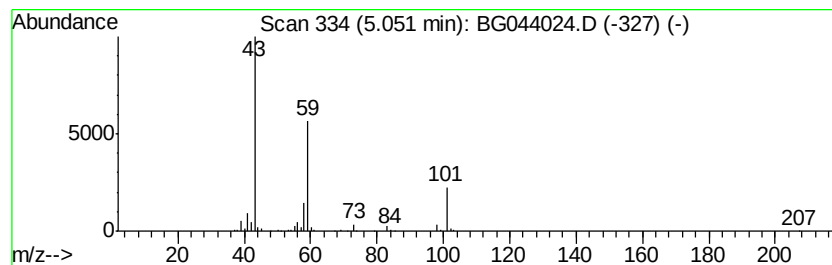
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ClientSampleId :
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.05	8.74 ng	278477	1,4-Dichlorobenzene-d4	7.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	50
3	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	43
4	2-Butanol, 1-methoxy-	104	C5H12O2	053778-73-7	25
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25



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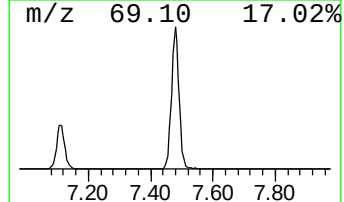
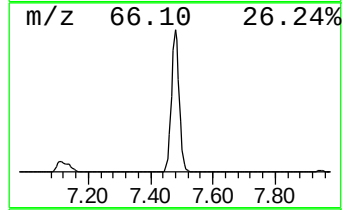
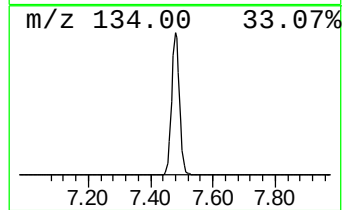
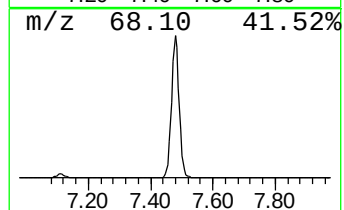
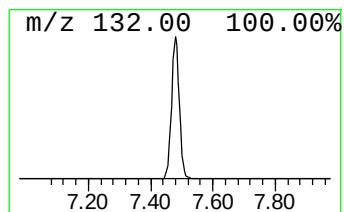
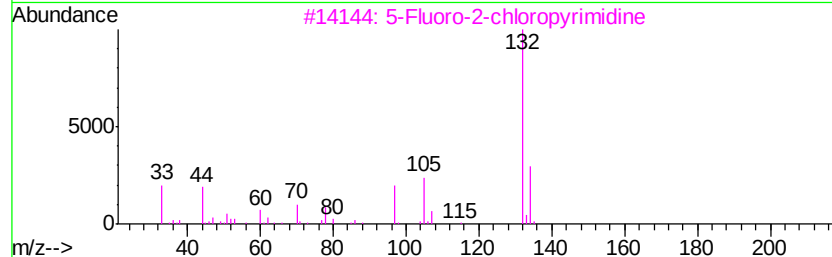
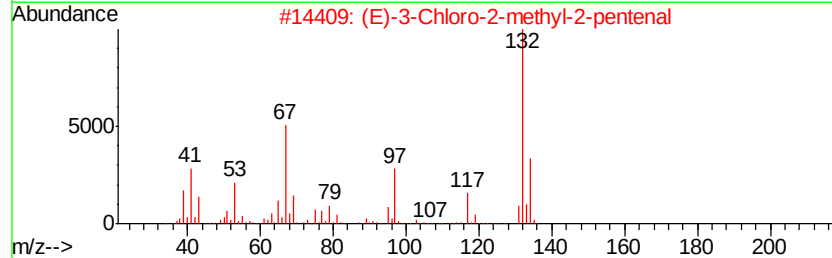
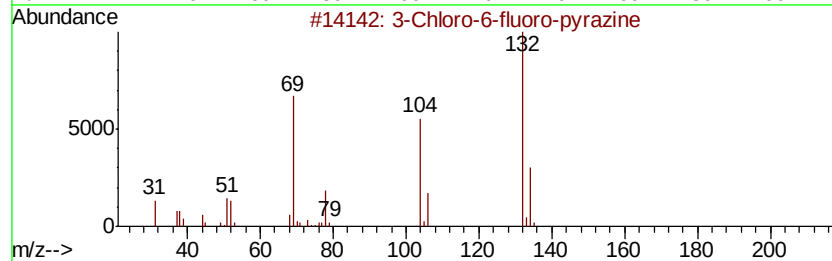
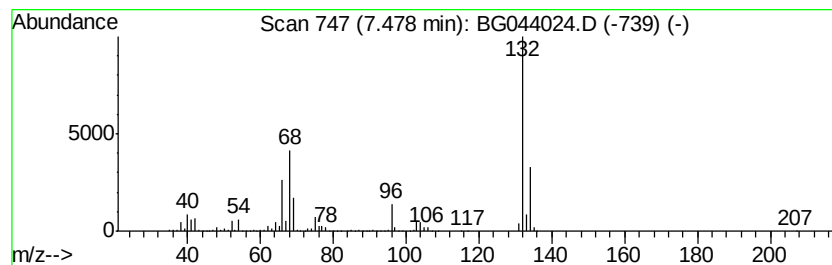
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Peak Number 3 unknown7.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	88.22 ng	2811620	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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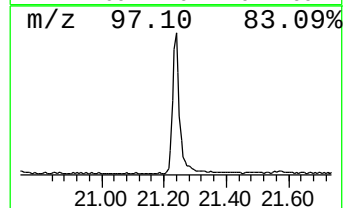
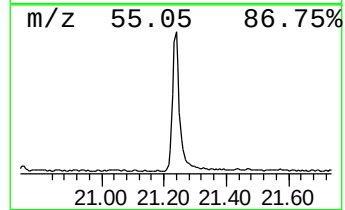
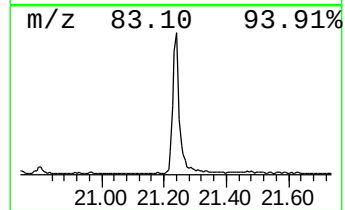
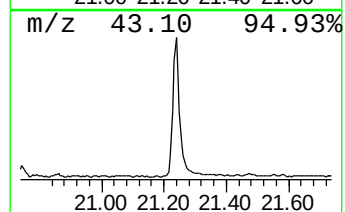
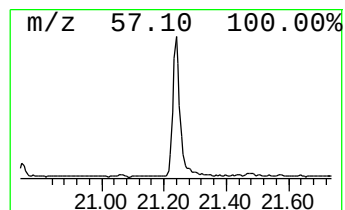
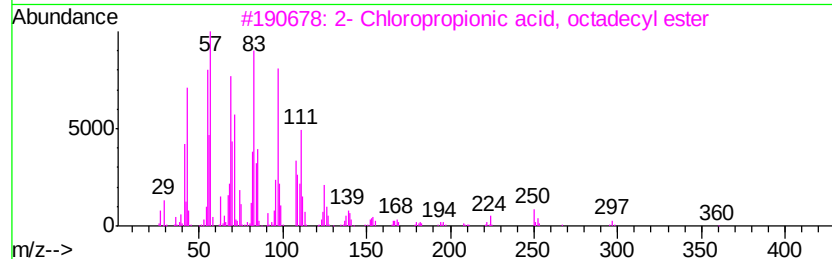
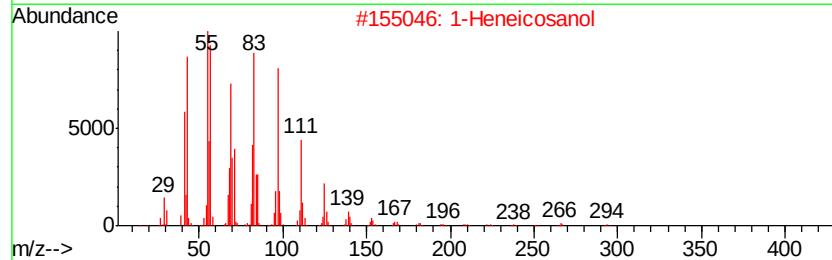
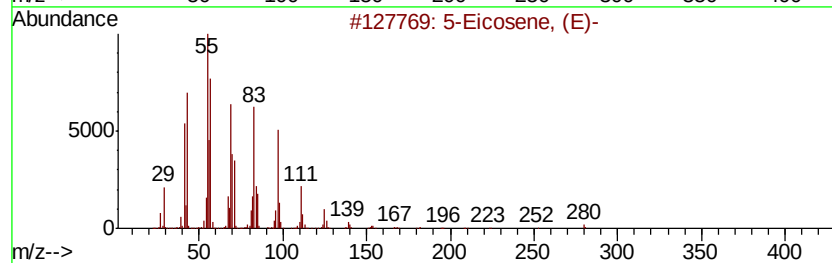
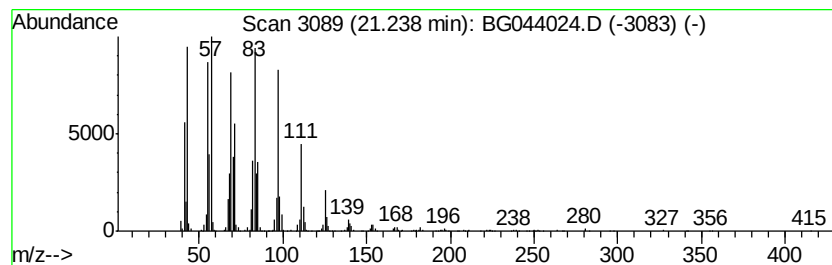
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Peak Number 5 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.24	15.13 ng	1246090	Chrysene-d12	21.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	98
2	1-Heneicosanol		312	C21H44O	015594-90-8	95
3	2- Chloropropionic acid, octadec...		360	C21H41ClO2	088104-31-8	94
4	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	94
5	1-Heneicosyl formate		340	C22H44O2	077899-03-7	94



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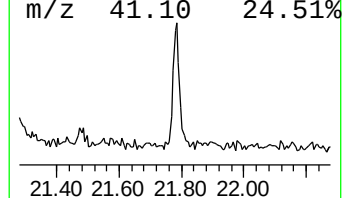
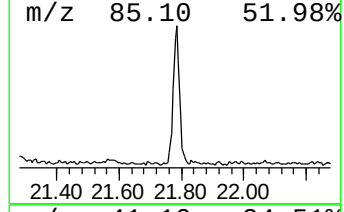
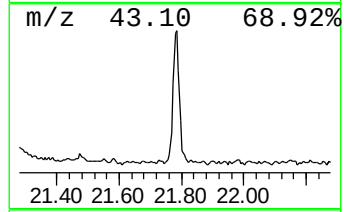
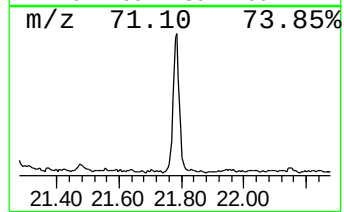
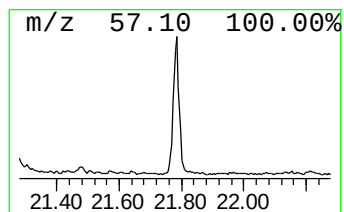
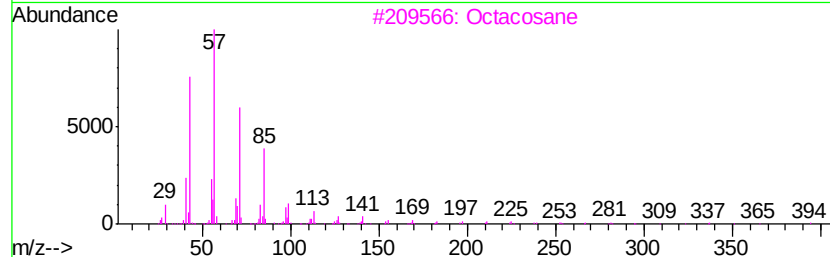
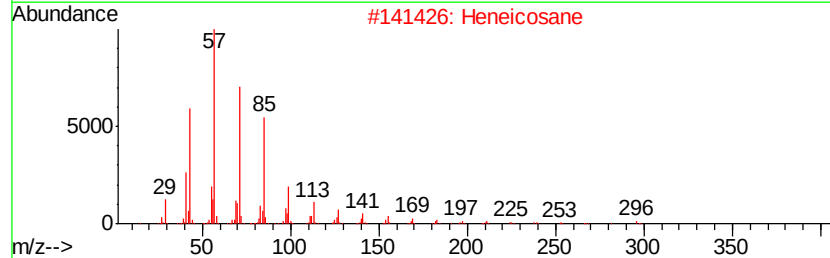
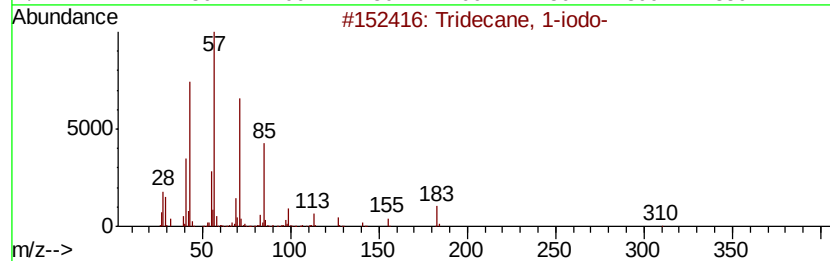
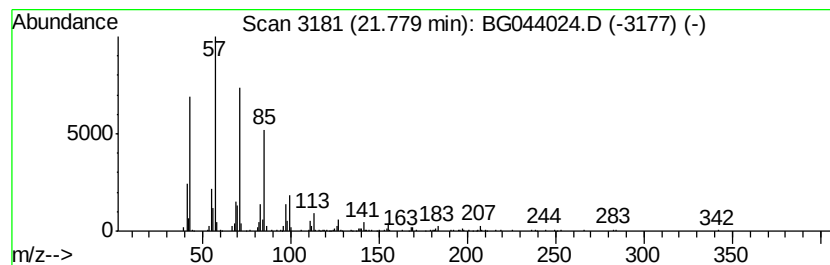
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Peak Number 6 Tridecane, 1-iodo- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.78	2.99 ng	245875	Chrysene-d12	21.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Hexatriacontane	507	C36H74	000630-06-8	91



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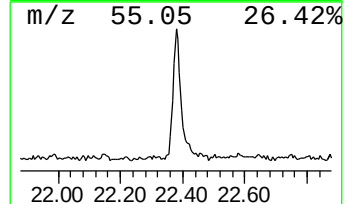
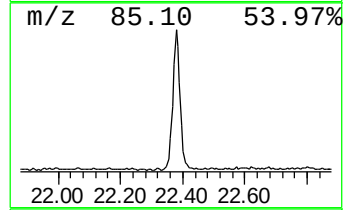
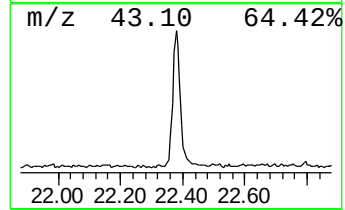
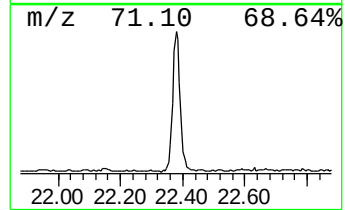
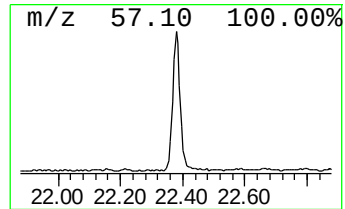
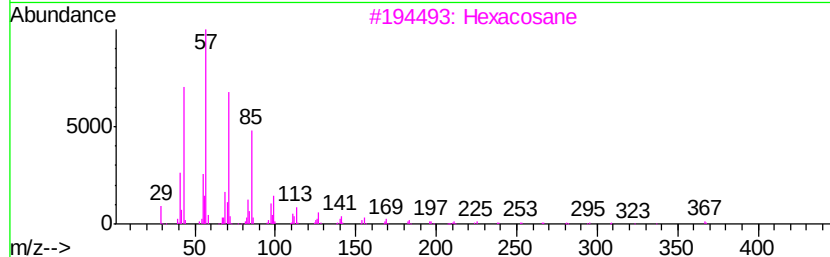
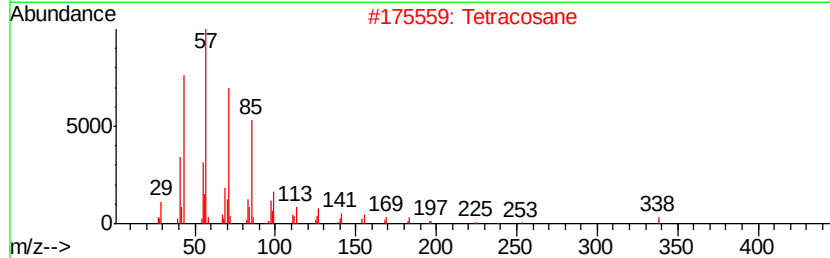
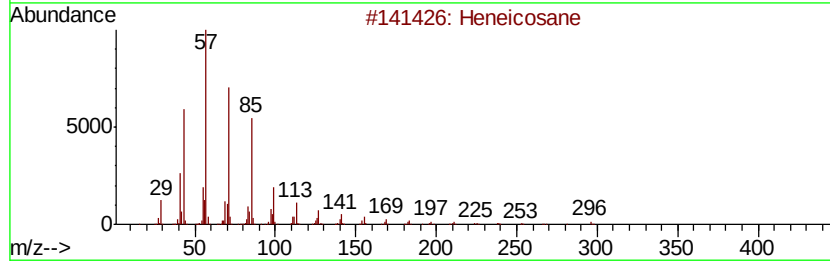
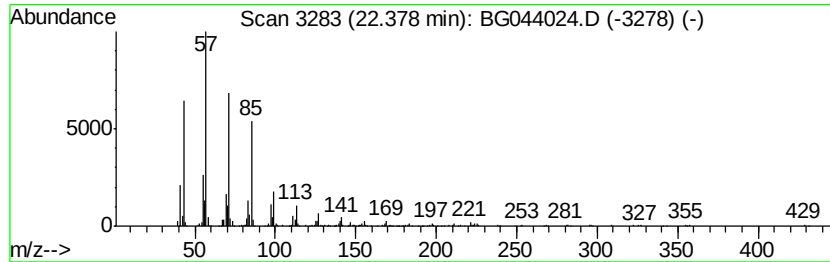
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.38	6.88 ng	566681	Chrysene-d12	21.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	97
2	Tetracosane		338	C24H50	000646-31-1	91
3	Hexacosane		366	C26H54	000630-01-3	91
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
5	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
 Data File : BG044024.D
 Acq On : 13 Jan 2020 22:57
 Operator : JU
 Sample : L1074-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 CIY1-GW-102-13.5

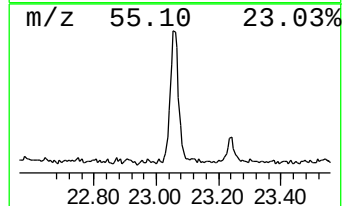
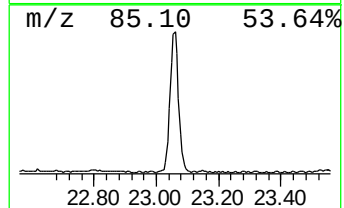
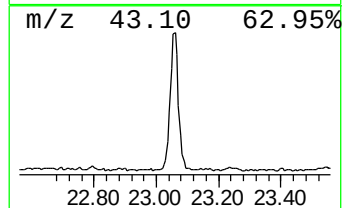
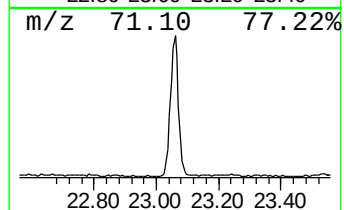
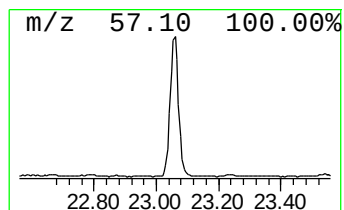
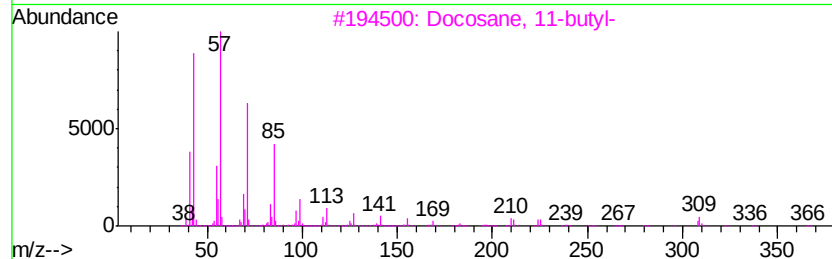
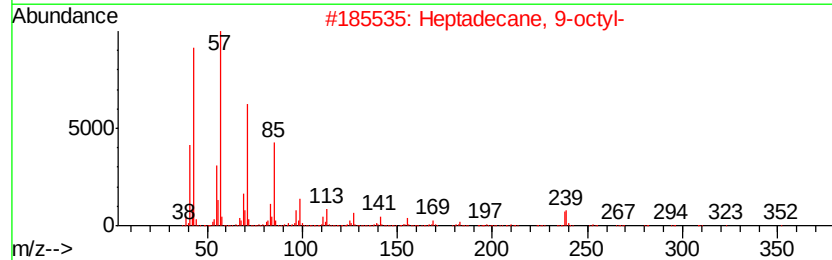
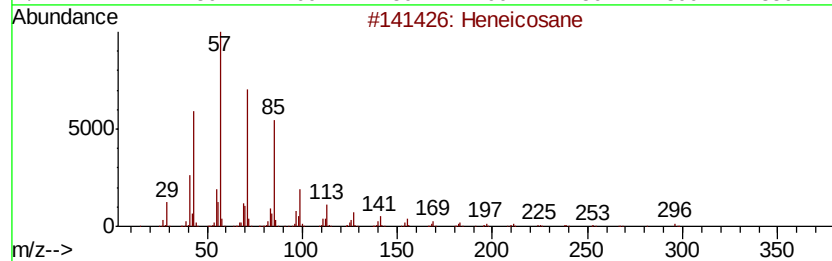
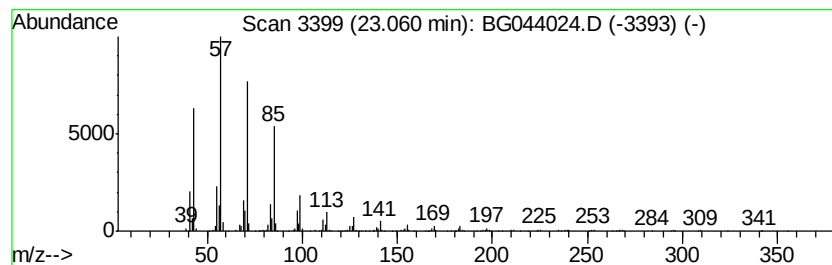
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Heptadecane, 9-octyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.06	8.03 ng	661071	Chrysene-d12	21.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
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Instrument :
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ClientSampleId :
CIY1-GW-102-13.5

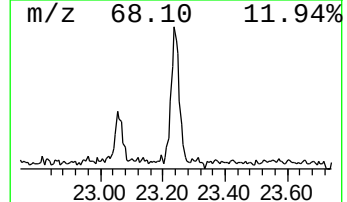
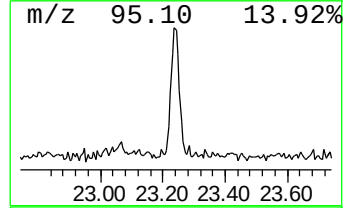
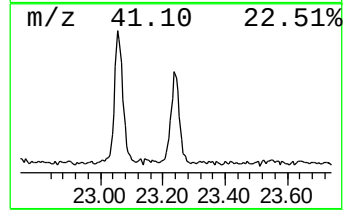
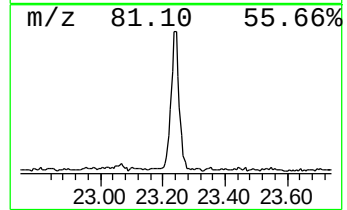
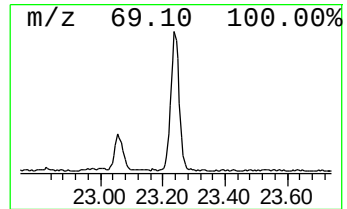
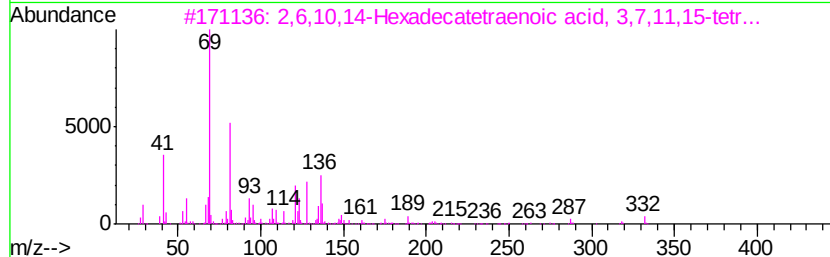
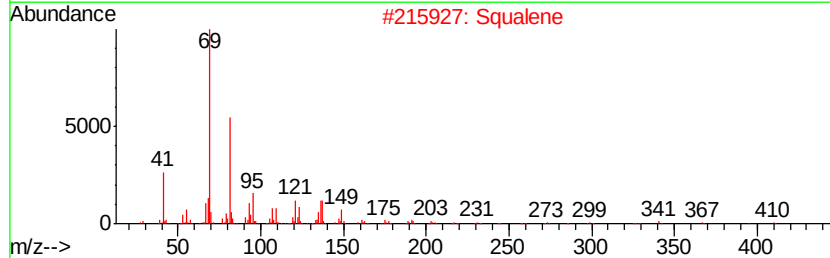
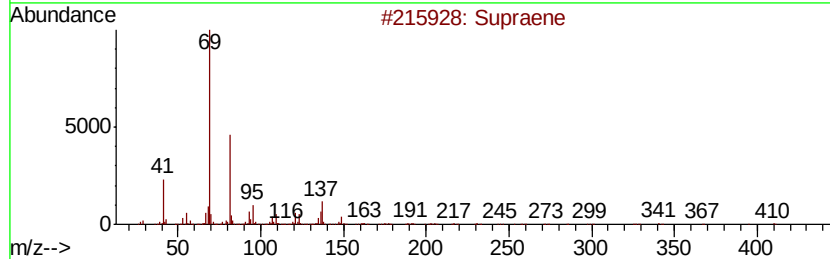
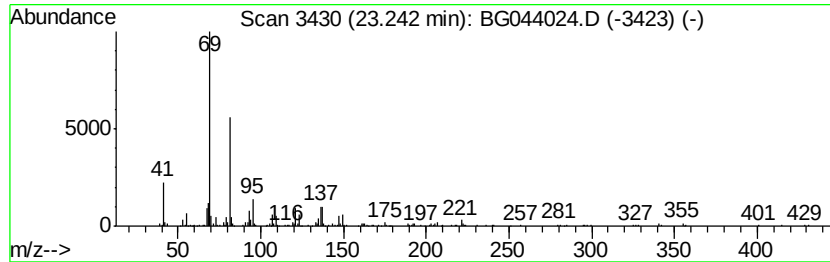
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Supraene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.24	4.15 ng	337431	Perylene-d12	24.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	91
3		2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	78
4		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	72
5		Farnesol isomer a	222	C15H26O	1000108-92-4	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
Data File : BG044024.D
Acq On : 13 Jan 2020 22:57
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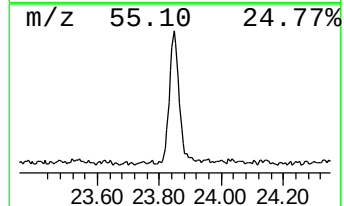
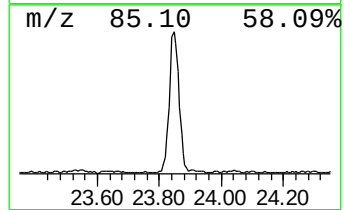
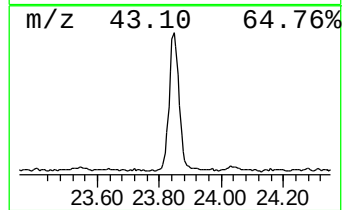
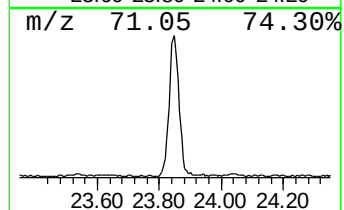
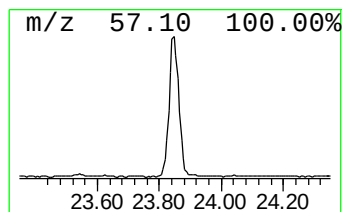
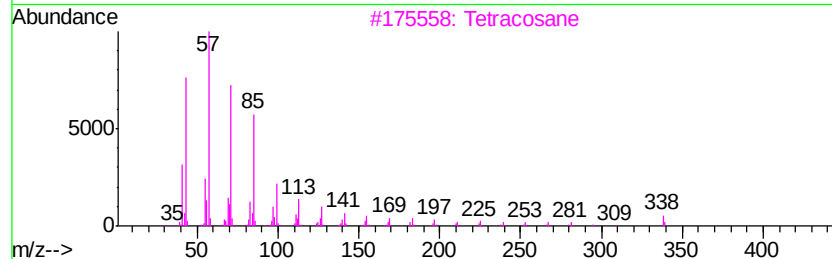
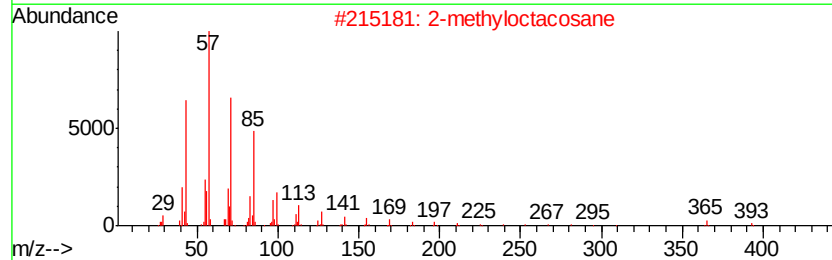
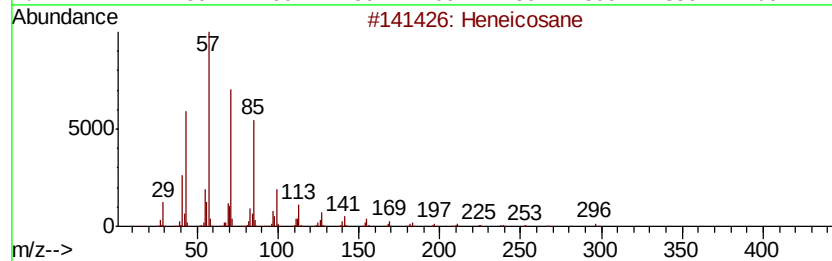
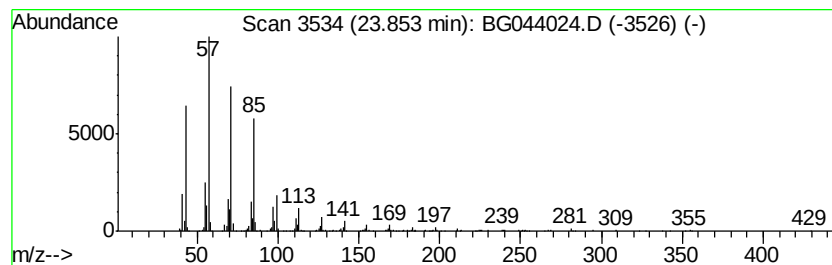
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-methyloctacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.85	9.68 ng	786334	Perylene-d12	24.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			2-methyloctacosane	408	C29H60	1000376-72-8	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5			Octadecane	254	C18H38	000593-45-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
Data File : BG044024.D
Acq On : 13 Jan 2020 22:57
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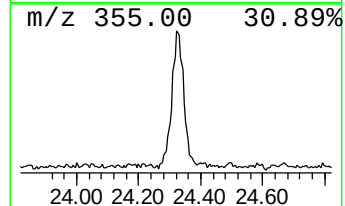
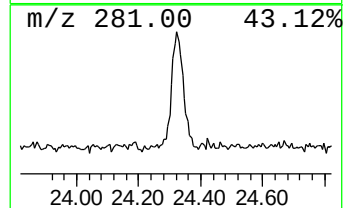
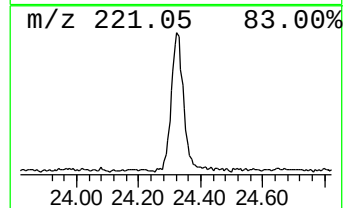
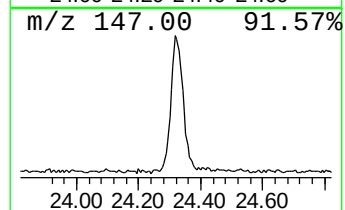
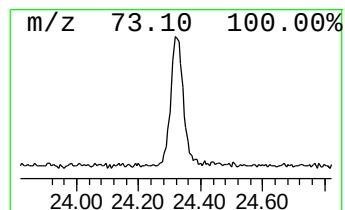
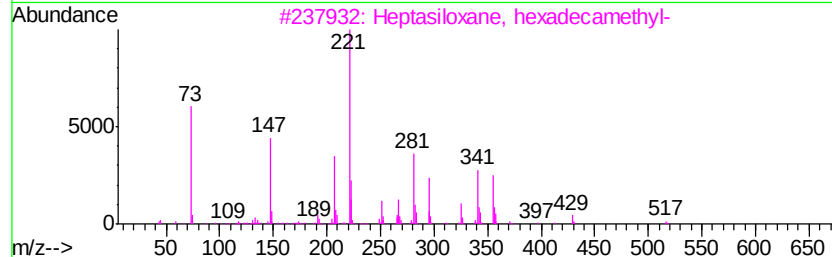
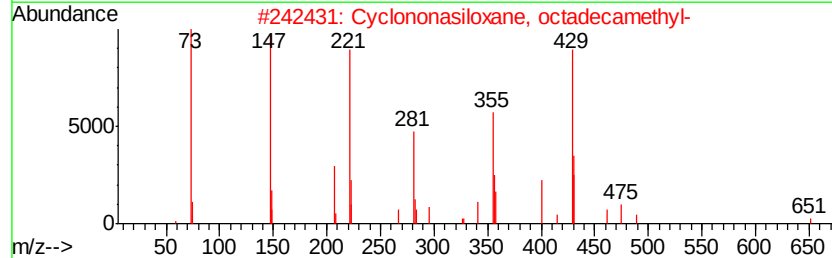
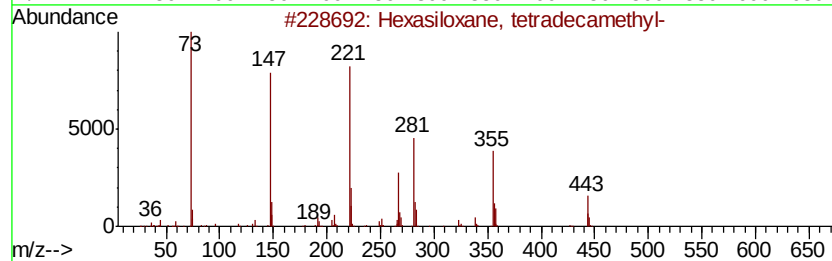
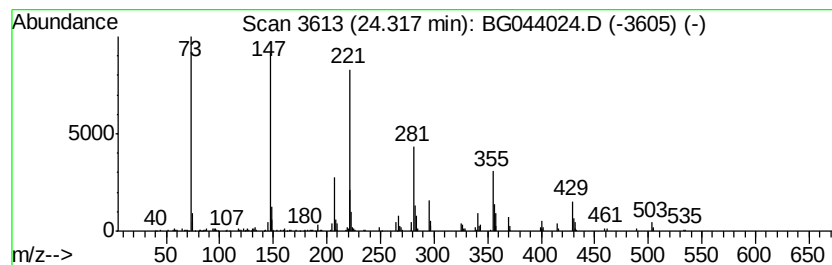
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexasiloxane, tetradecamethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.32	5.28 ng	429154	Perylene-d12	24.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	72
2		Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	53
3		Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	43
4		1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	37
5		3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
 Data File : BG044024.D
 Acq On : 13 Jan 2020 22:57
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Instrument :
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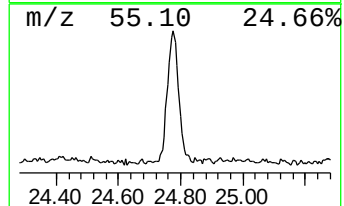
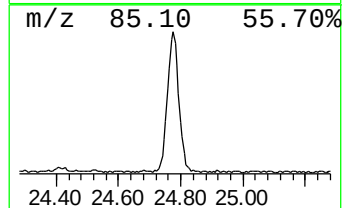
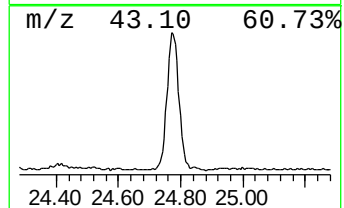
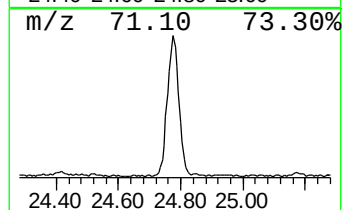
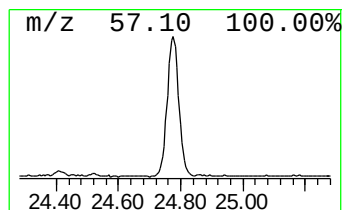
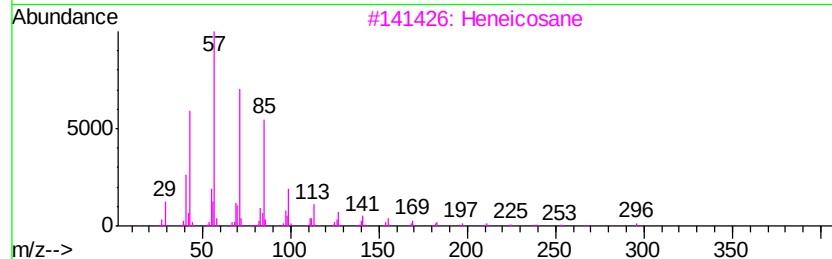
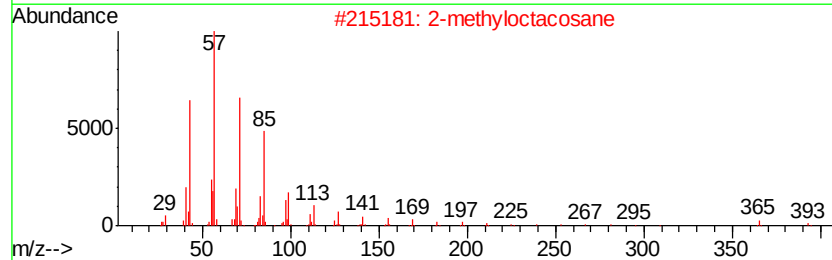
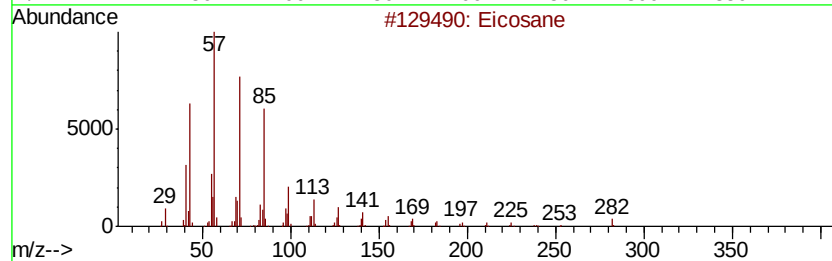
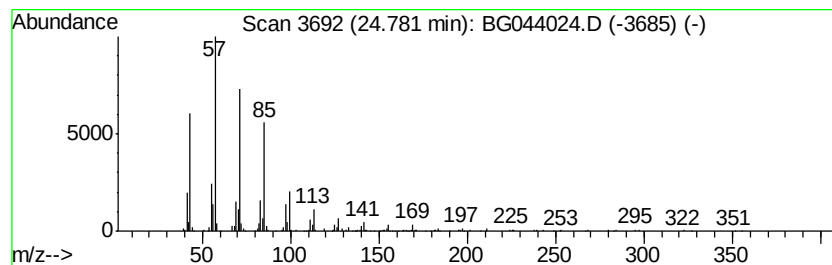
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Eicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.78	11.03 ng	896461	Perylene-d12	24.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Octadecane	254	C18H38	000593-45-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
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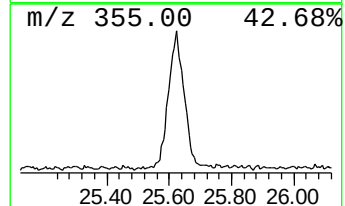
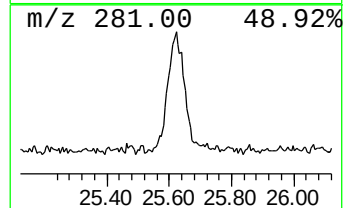
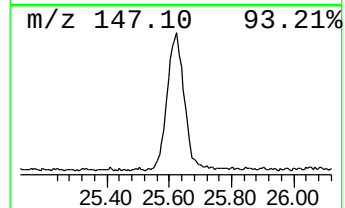
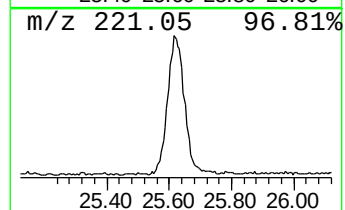
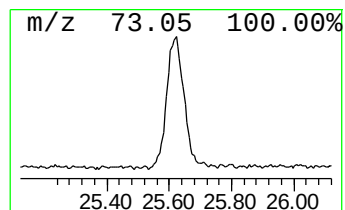
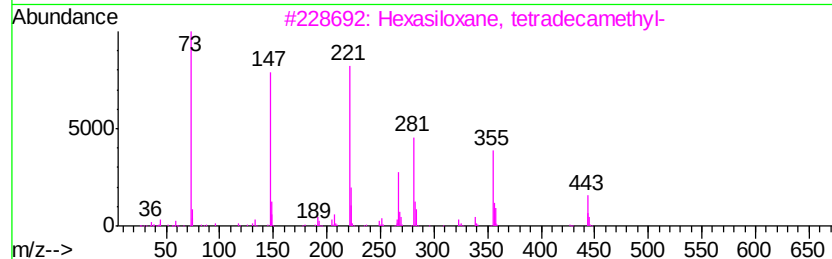
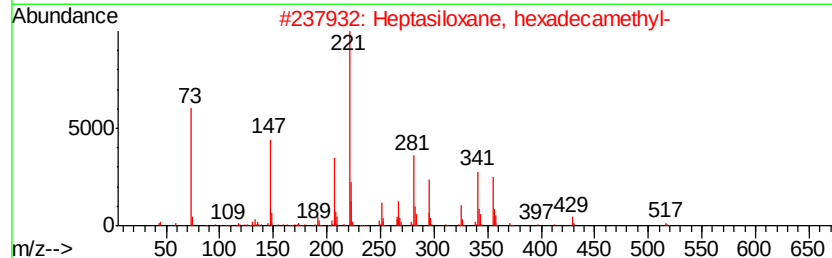
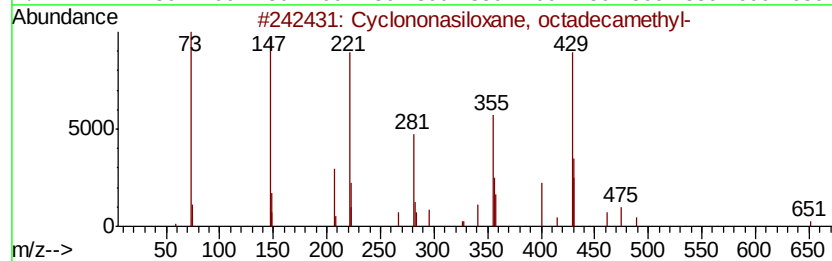
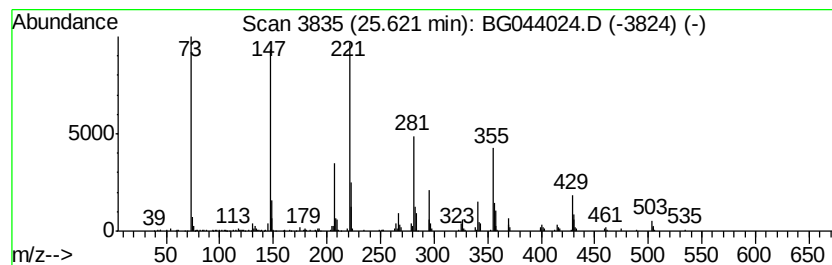
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclononasiloxane, octadeca... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.62	8.69 ng	705846	Perylene-d12	24.69	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	87
2	Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	86
3	Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	72
4	3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	47
5	3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
Data File : BG044024.D
Acq On : 13 Jan 2020 22:57
Operator : JU
Sample : L1074-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CIY1-GW-102-13.5

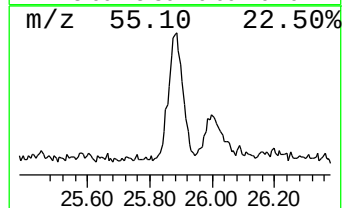
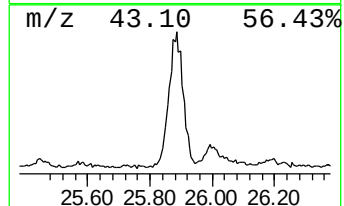
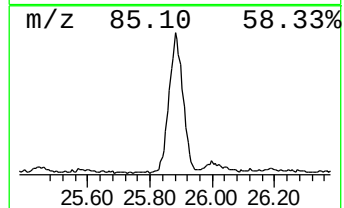
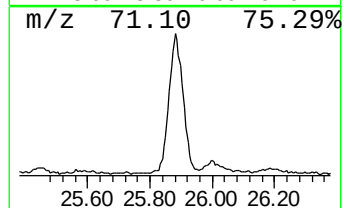
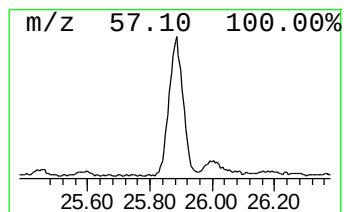
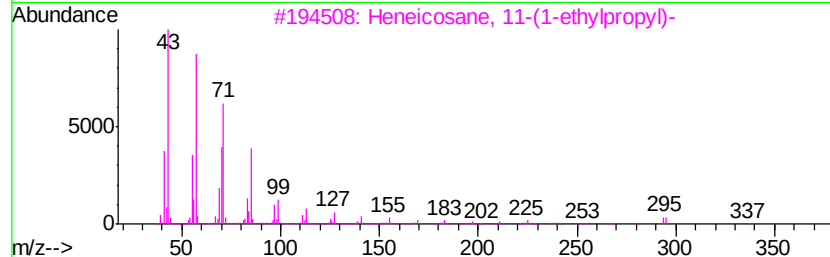
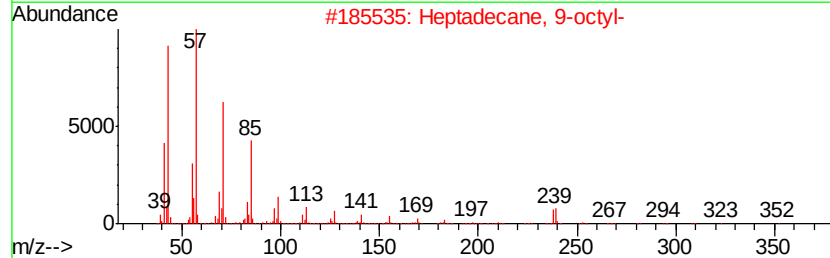
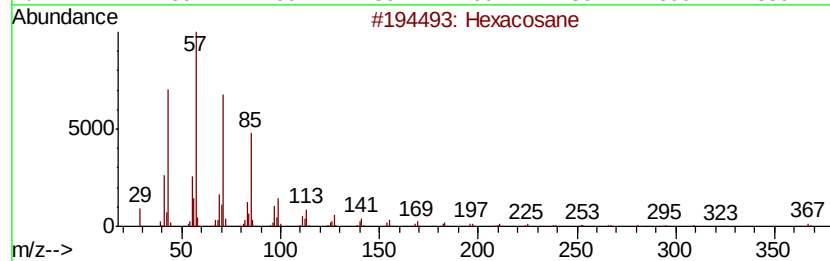
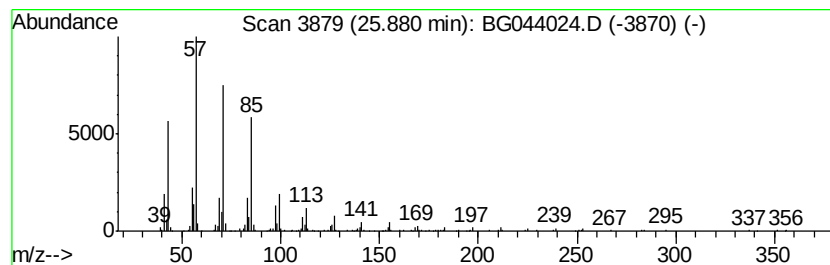
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.88	8.29 ng	673114	Perylene-d12	24.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
 Data File : BG044024.D
 Acq On : 13 Jan 2020 22:57
 Operator : JU
 Sample : L1074-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

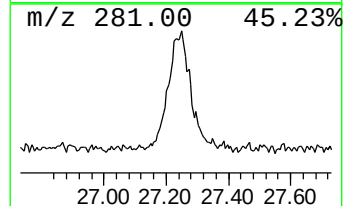
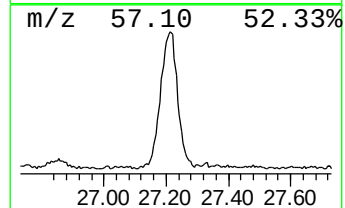
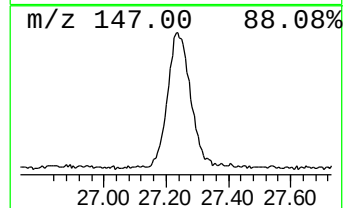
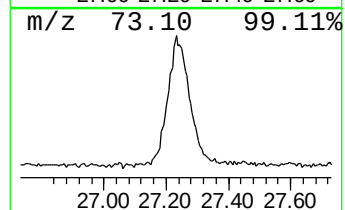
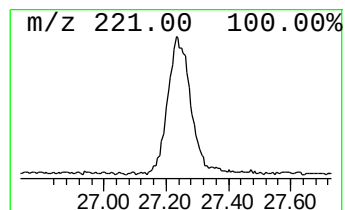
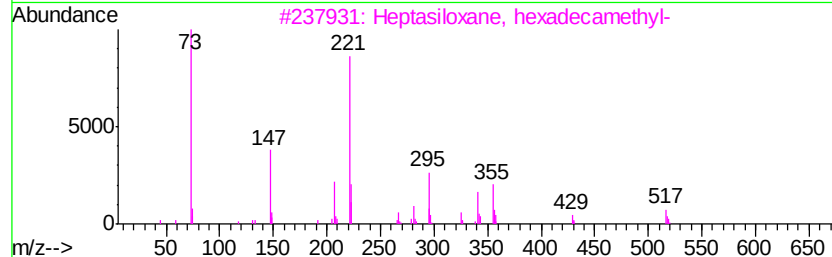
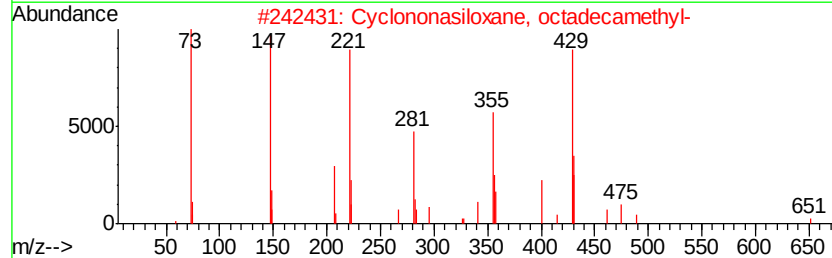
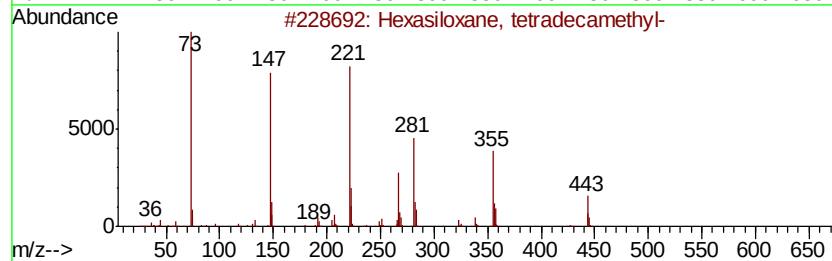
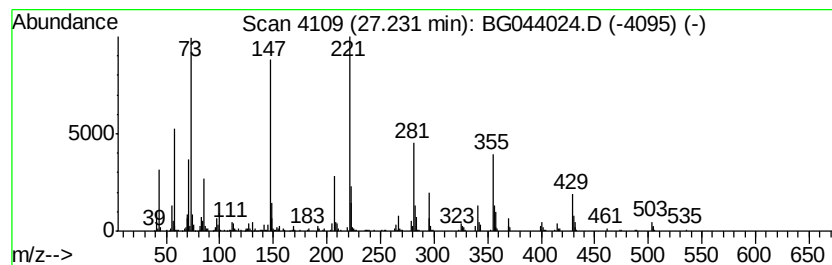
Instrument :
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 ClientSampleId :
 CIY1-GW-102-13.5

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

 Peak Number 15 Hexasiloxane, tetradecamethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
27.23	16.59 ng	1347880	Perylene-d12	24.69	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	76
2	Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	50
3	Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	35
4	N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	25
5	Dithioerythritol, 0,0',S,S'-tetr...	442	C16H42O2S2Si4	1000079-30-7	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011320\
Data File : BG044024.D
Acq On : 13 Jan 2020 22:57
Operator : JU
Sample : L1074-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CIY1-GW-102-13.5

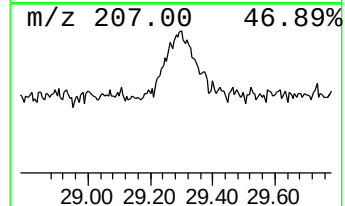
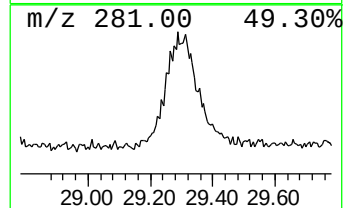
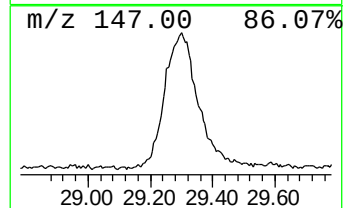
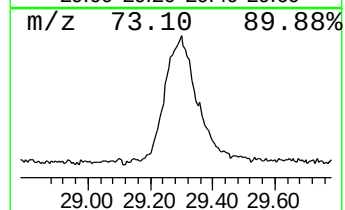
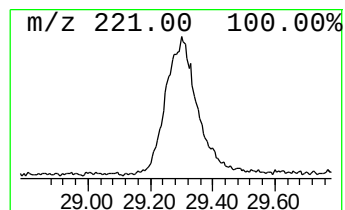
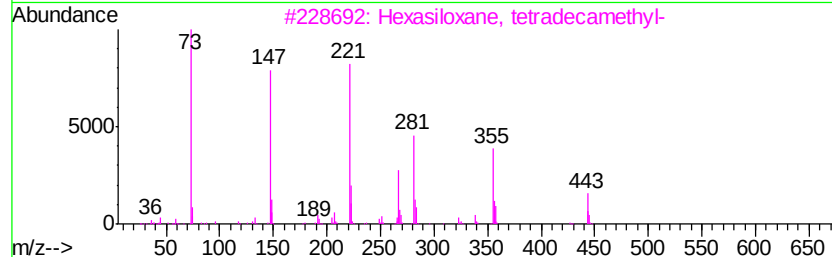
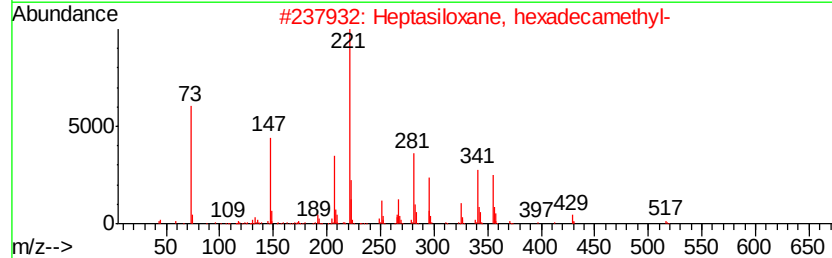
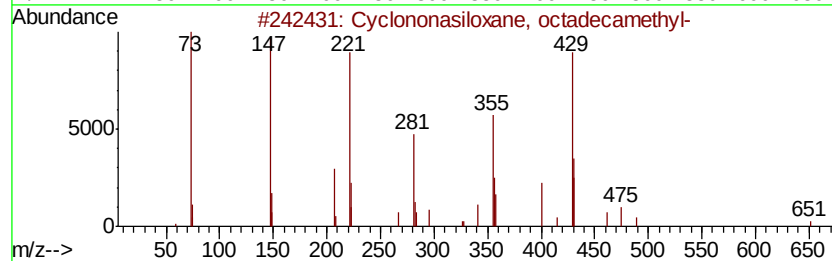
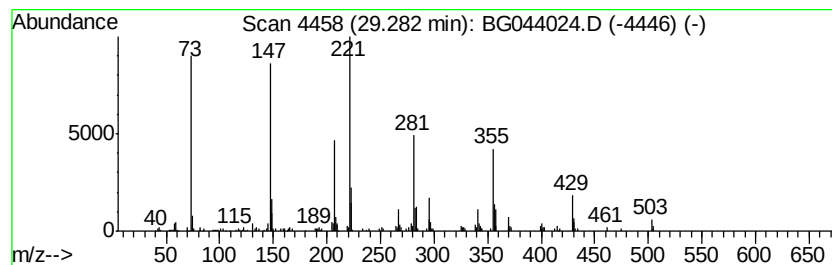
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Cyclononasiloxane, octadeca... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.28	3.46 ng	281025	Perylene-d12	24.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H5409Si9	000556-71-8	83
2			Heptasiloxane, hexadecamethyl-	532	C16H4806Si7	000541-01-5	78
3			Hexasiloxane, tetradecamethyl-	458	C14H4205Si6	000107-52-8	64
4			Pentasiloxane, dodecamethyl-	384	C12H3604Si5	000141-63-9	22
5			Silane, [[(3.beta.,11.beta.,20S)...]	638	C33H6604Si4	032221-72-0	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG011320\
 Data File : BG044024.D
 Acq On : 13 Jan 2020 22:57
 Operator : JU
 Sample : L1074-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CIY1-GW-102-13.5

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...	3.12	77.9	ng	2481480	1	7.95	637383	20.0
2-Pentanone, 4-hy...	5.05	8.7	ng	278477	1	7.95	637383	20.0
unknown7.48	7.48	88.2	ng	2811620	1	7.95	637383	20.0
5-Eicosene, (E)-	21.24	15.1	ng	1246090	5	21.57	1647290	20.0
Tridecane, 1-iodo-	21.78	3.0	ng	245875	5	21.57	1647290	20.0
Heneicosane	22.38	6.9	ng	566681	5	21.57	1647290	20.0
Heptadecane, 9-oc...	23.06	8.0	ng	661071	5	21.57	1647290	20.0
Supraene	23.24	4.2	ng	337431	6	24.69	1624790	20.0
2-methyloctacosane	23.85	9.7	ng	786334	6	24.69	1624790	20.0
Hexasiloxane, tet...	24.32	5.3	ng	429154	6	24.69	1624790	20.0
Eicosane	24.78	11.0	ng	896461	6	24.69	1624790	20.0
Cyclononasiloxane...	25.62	8.7	ng	705846	6	24.69	1624790	20.0
Hexacosane	25.88	8.3	ng	673114	6	24.69	1624790	20.0
Hexasiloxane, tet...	27.23	16.6	ng	1347880	6	24.69	1624790	20.0
Cyclononasiloxane...	29.28	3.5	ng	281025	6	24.69	1624790	20.0