

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
 Data File : BG044108.D
 Acq On : 20 Jan 2020 21:38
 Operator : JU
 Sample : L1165-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 QVD-GW-01-45

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	5.035	325	331	343	rVB	106736	172018	3.08%	0.575%
2	5.511	405	412	424	rBV	569669	961290	17.23%	3.214%
3	6.011	491	497	505	rBV	79022	130828	2.34%	0.437%
4	7.097	675	682	697	rBV	389977	722995	12.96%	2.417%
5	7.468	737	745	757	rBV	1146767	2009875	36.02%	6.720%
6	7.932	816	824	831	rBV	267713	448308	8.03%	1.499%
7	8.255	870	879	887	rBV	1058178	1840342	32.98%	6.154%
8	9.083	1010	1020	1031	rBV	830301	1542020	27.63%	5.156%
9	10.734	1292	1301	1315	rBV	380101	689872	12.36%	2.307%
10	13.190	1710	1719	1733	rBV	2454561	3875832	69.45%	12.960%
11	13.472	1761	1767	1775	rVB2	71405	116795	2.09%	0.391%
12	14.019	1854	1860	1871	rBV	135389	213793	3.83%	0.715%
13	14.565	1943	1953	1963	rBV	659595	1070177	19.18%	3.578%
14	16.058	2199	2207	2219	rBV	1981077	3192829	57.21%	10.676%
15	16.152	2219	2223	2236	rVB2	25293	64941	1.16%	0.217%
16	17.309	2413	2420	2431	rBV	824783	1282768	22.99%	4.289%
17	19.072	2716	2720	2728	rBV2	46955	94548	1.69%	0.316%
18	19.389	2765	2774	2778	rBV3	35882	59675	1.07%	0.200%
19	19.924	2859	2865	2879	rBV	3986467	5580594	100.00%	18.660%
20	20.247	2912	2920	2924	rBV	48578	79747	1.43%	0.267%
21	20.740	3000	3004	3008	rVB	48439	57587	1.03%	0.193%
22	21.234	3082	3088	3101	rBV	240386	502377	9.00%	1.680%
23	21.557	3136	3143	3152	rBV	794702	1331808	23.86%	4.453%
24	21.769	3174	3179	3185	rBV	131334	184955	3.31%	0.618%
25	22.362	3274	3280	3285	rBV	198237	318889	5.71%	1.066%
26	23.032	3388	3394	3404	rVB	233793	437436	7.84%	1.463%
27	23.813	3520	3527	3535	rBV	226775	475664	8.52%	1.590%
28	24.665	3661	3672	3679	rBV	483640	1384474	24.81%	4.629%
29	24.730	3679	3683	3693	rVB2	191773	463406	8.30%	1.549%
30	25.823	3860	3869	3880	rVB2	139728	395069	7.08%	1.321%
31	27.133	4083	4092	4102	rBV3	64155	206233	3.70%	0.690%

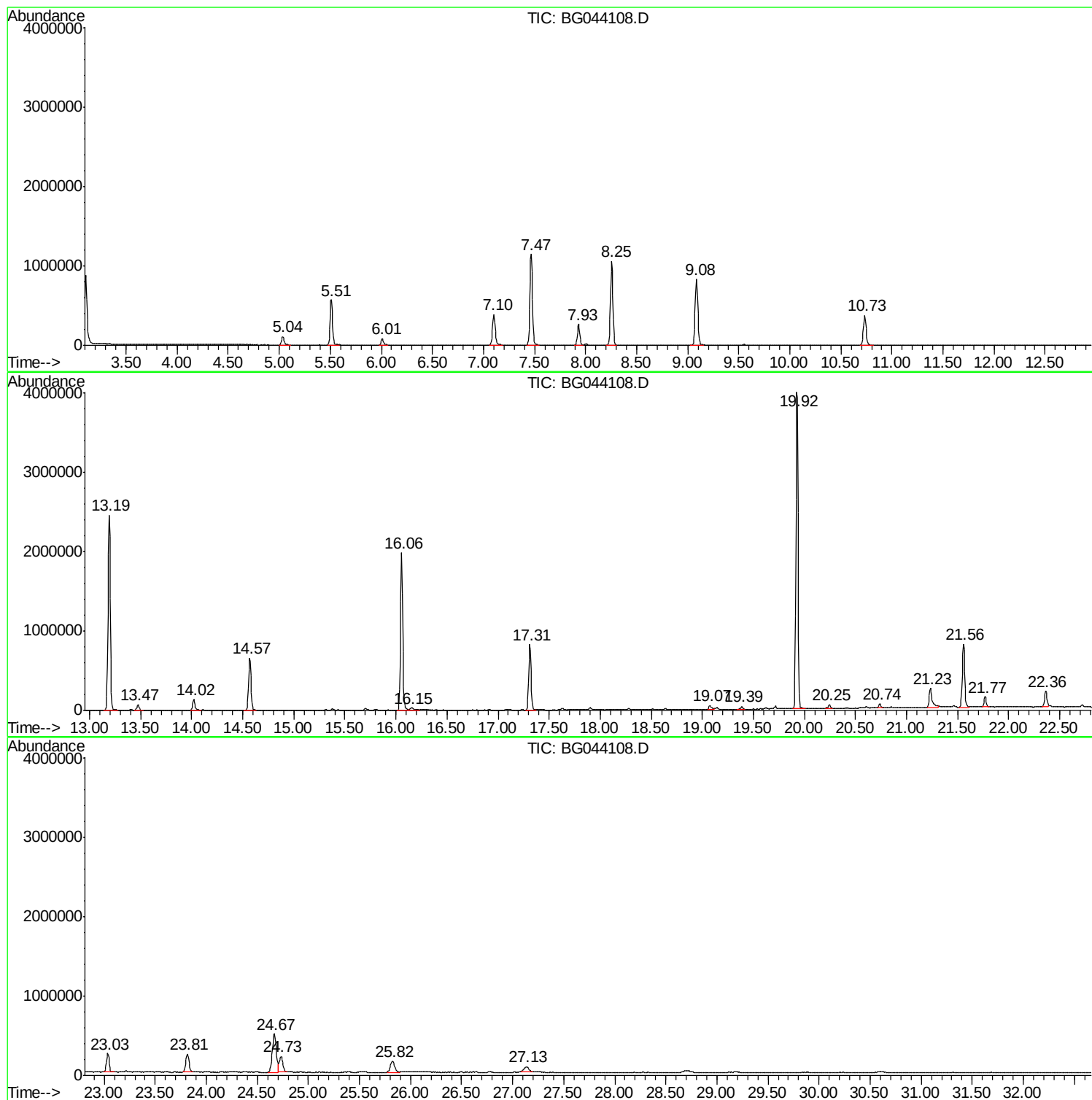
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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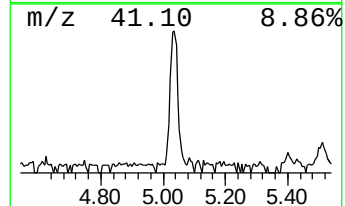
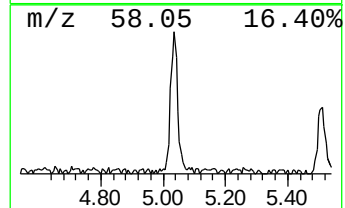
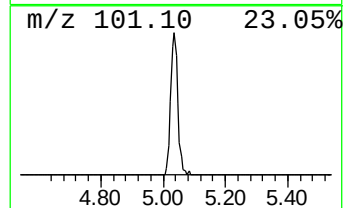
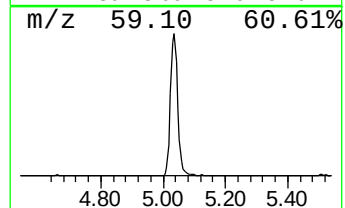
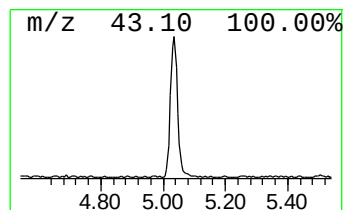
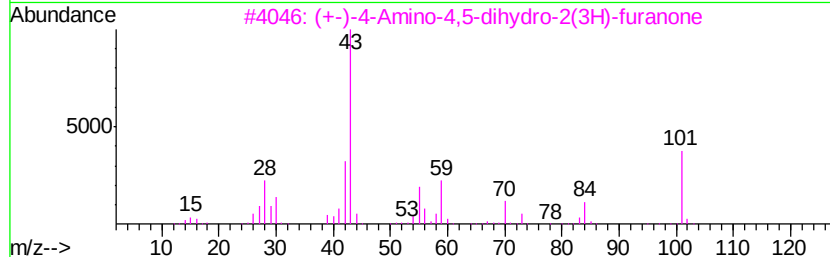
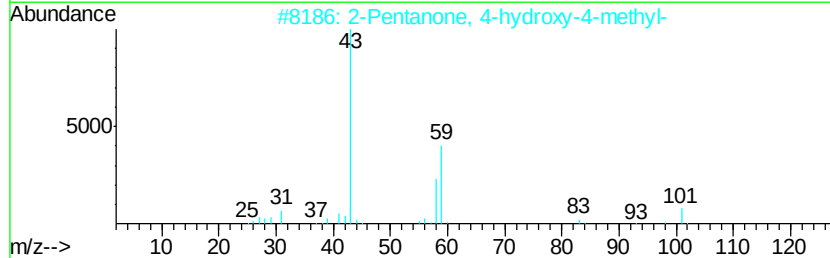
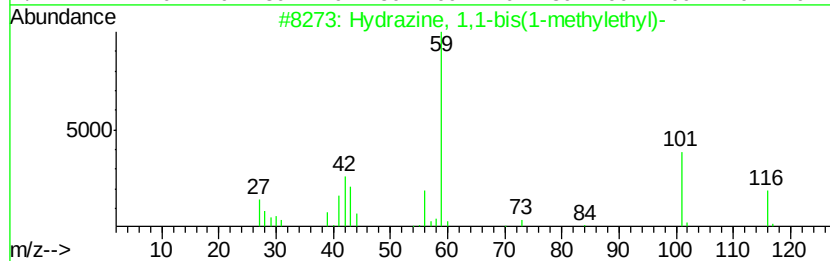
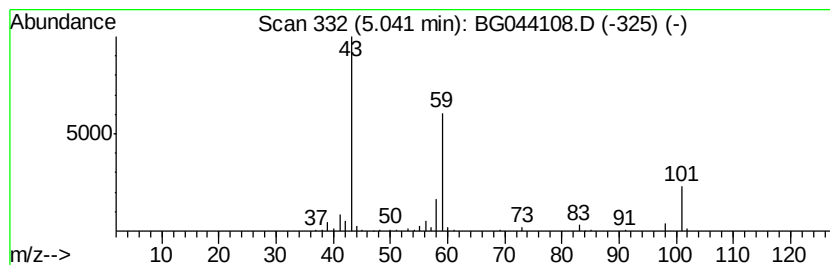
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown5.04 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	7.67 ng	172018	1,4-Dichlorobenzene-d4	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	40
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
3		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	32
4		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25



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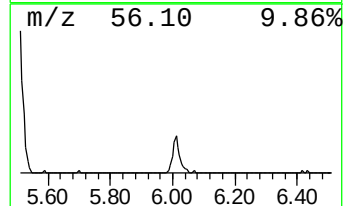
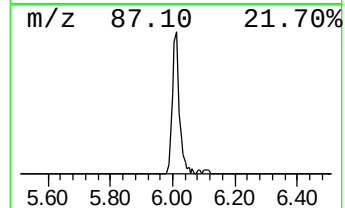
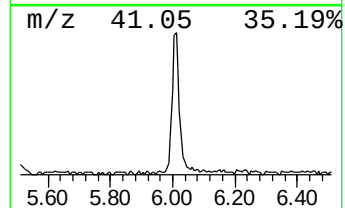
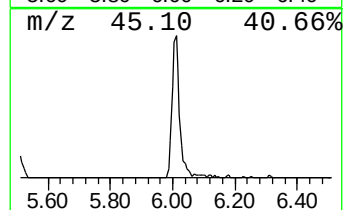
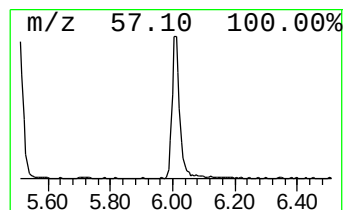
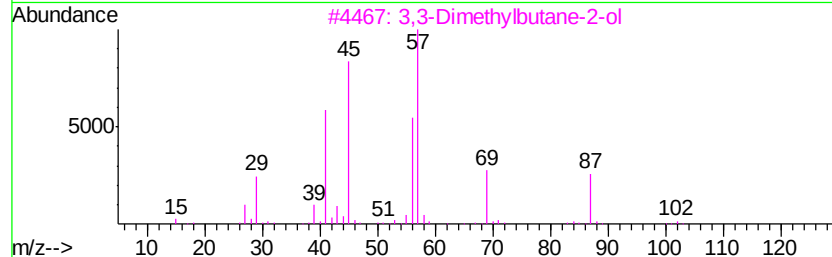
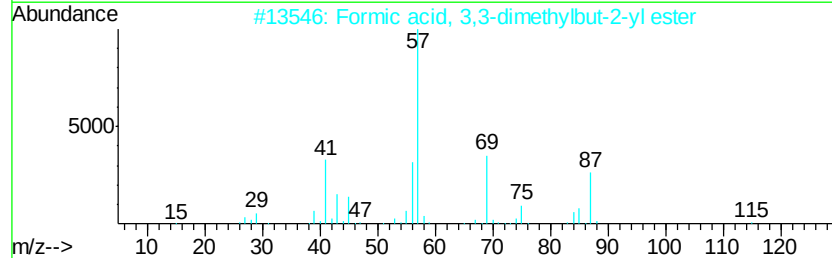
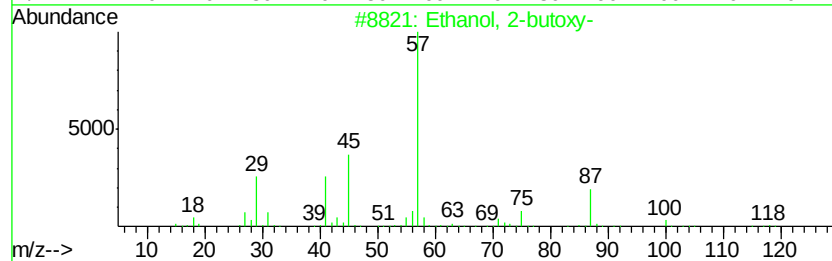
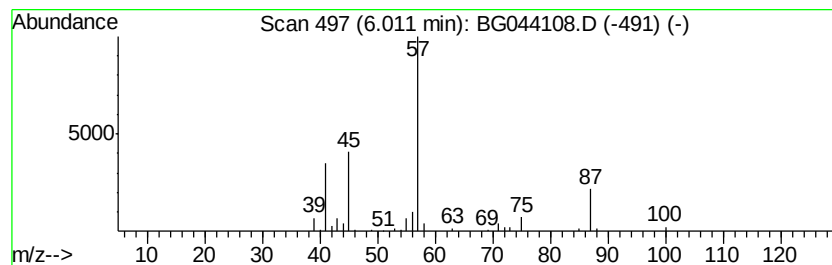
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TIC Library : C:\Database\NIST11.L
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Peak Number 2 Ethanol, 2-butoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.01	5.84 ng	130828	1,4-Dichlorobenzene-d4	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
2		Formic acid, 3,3-dimethylbut-2-yl...	130	C7H14O2	1000368-20-5	50
3		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	33
4		Propanoic acid, 2-methylpropyl e...	130	C7H14O2	000540-42-1	12
5		Propanoic acid, 2,2-dimethyl-	102	C5H10O2	000075-98-9	10



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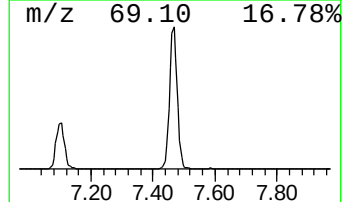
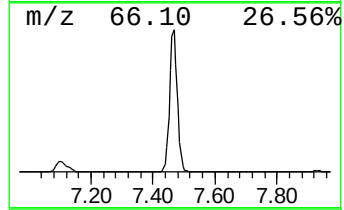
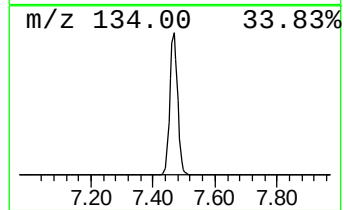
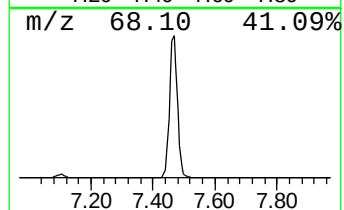
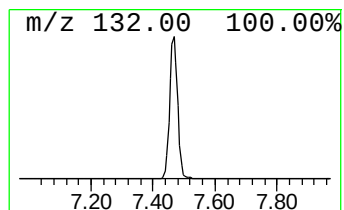
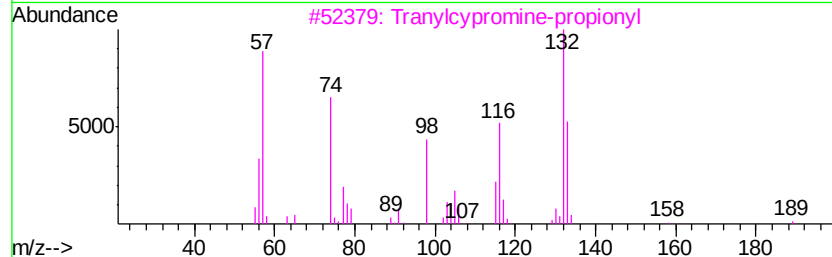
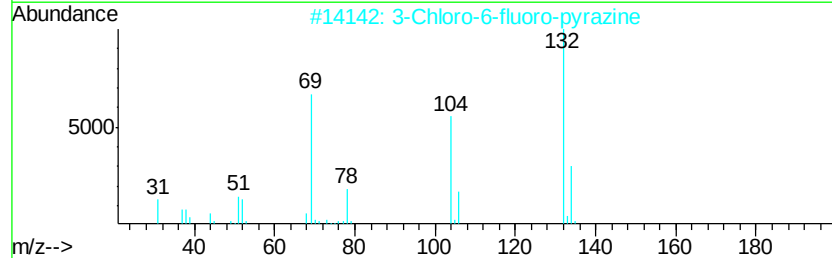
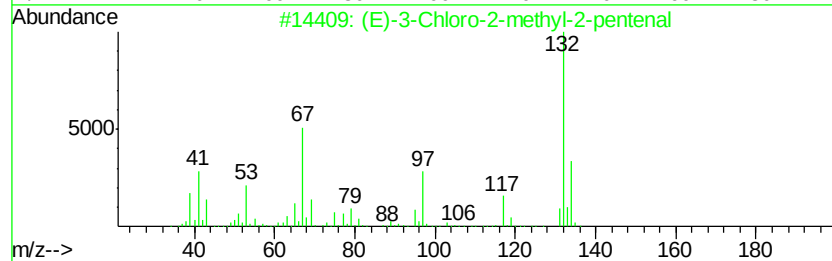
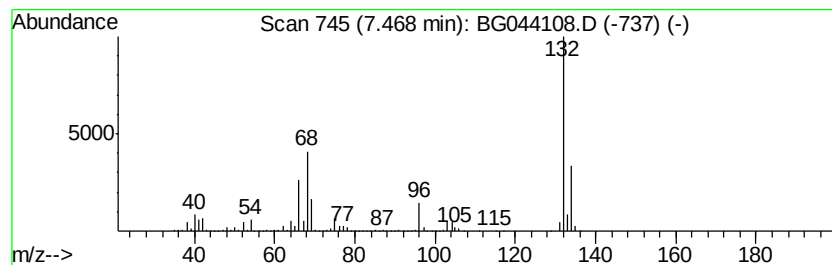
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Peak Number 3 unknown7.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	89.66 ng	2009880	1,4-Dichlorobenzene-d4	7.93

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



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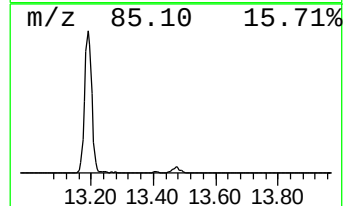
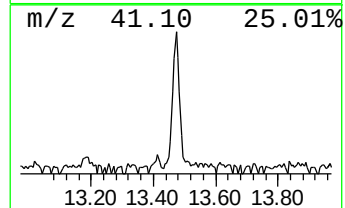
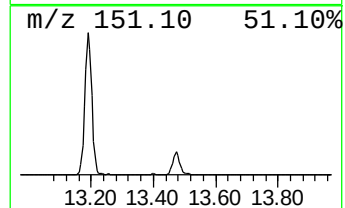
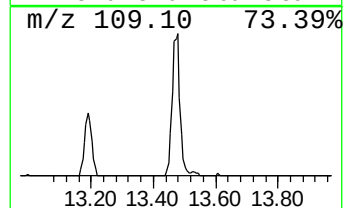
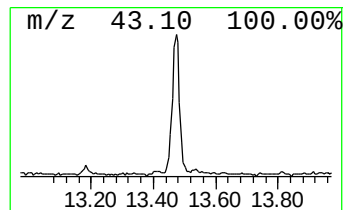
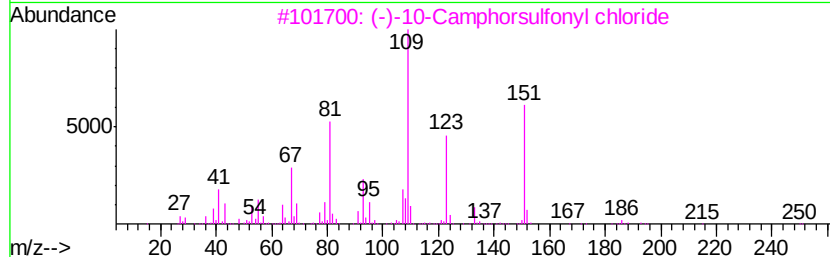
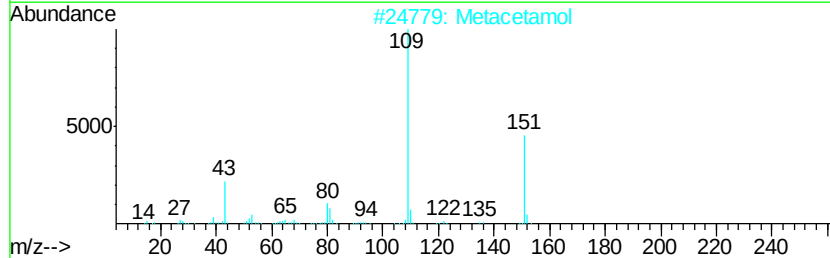
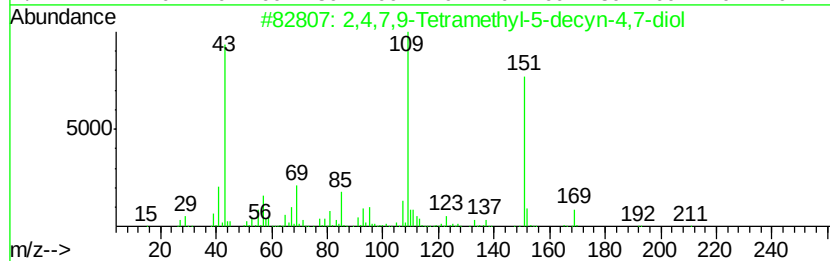
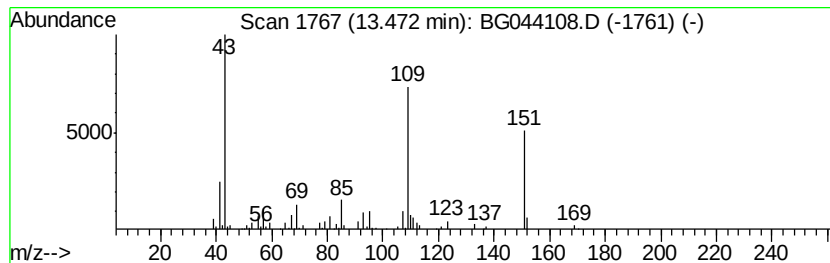
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Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.47	2.18 ng	116795	Acenaphthene-d10	14.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	Metacetamol	151	C8H9NO2	000621-42-1	58
3	(-)-10-Camphorsulfonyl chloride	250	C10H15ClO3S	039262-22-1	56
4	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	53
5	Acetaminophen	151	C8H9NO2	000103-90-2	42



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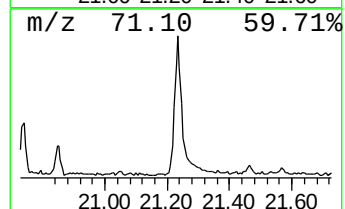
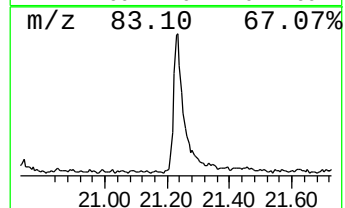
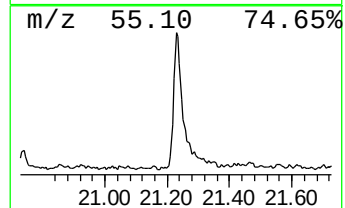
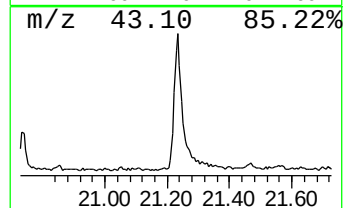
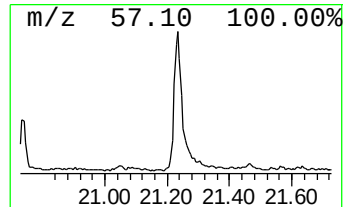
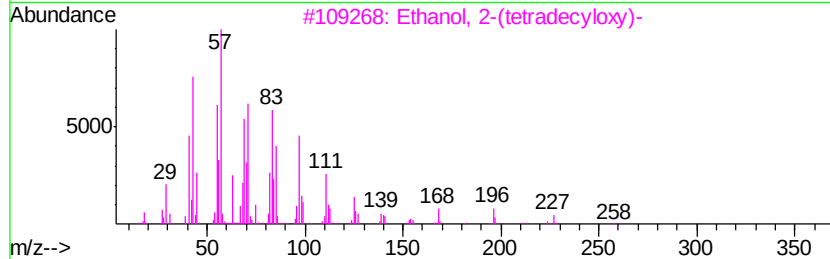
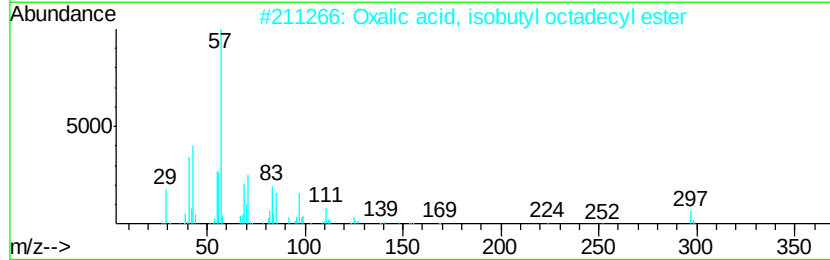
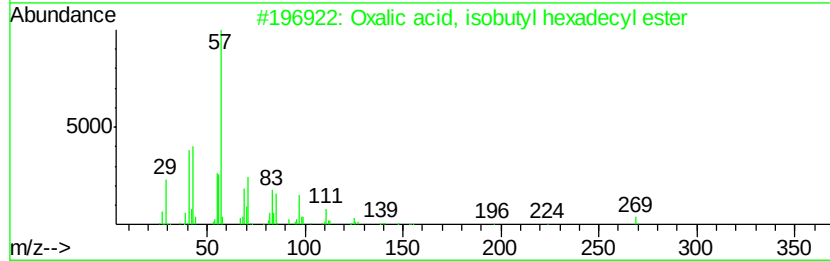
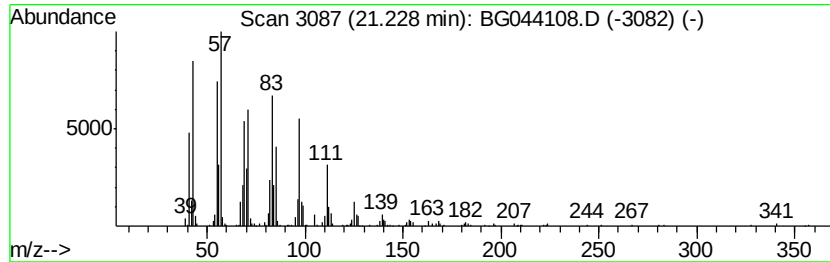
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Peak Number 6 Oxalic acid, isobutyl hexadecyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.23	7.54 ng	502377	Chrysene-d12	21.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
2		Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
3		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
4		Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	87
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	81



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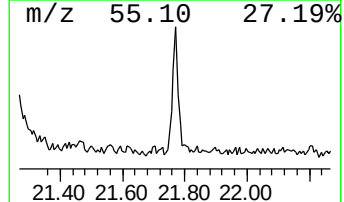
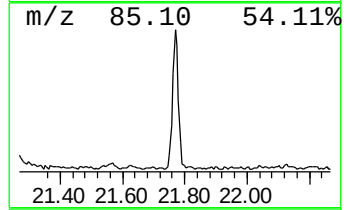
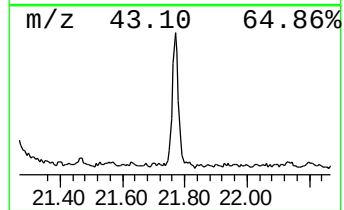
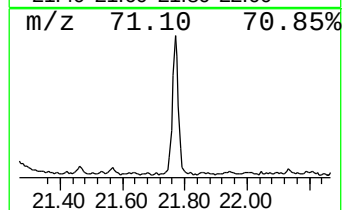
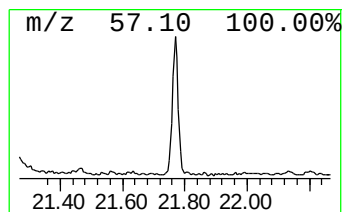
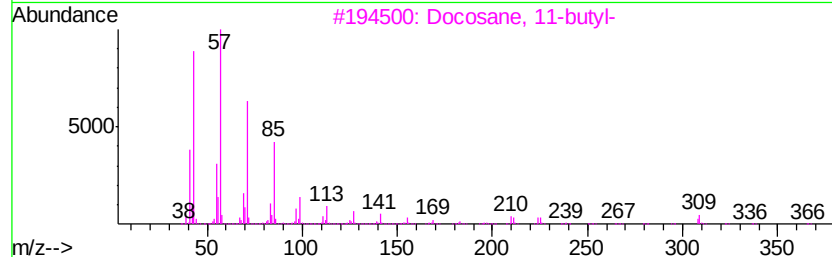
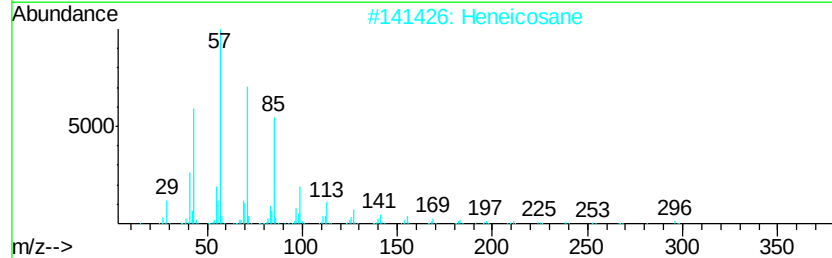
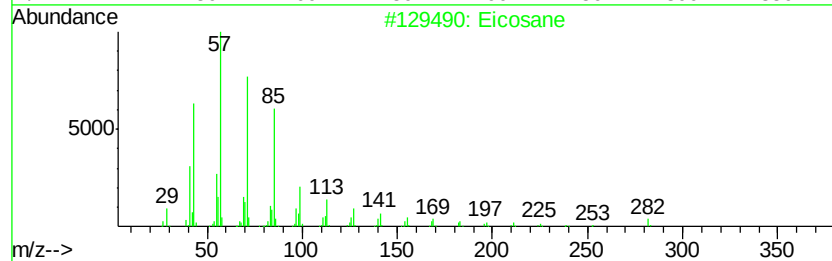
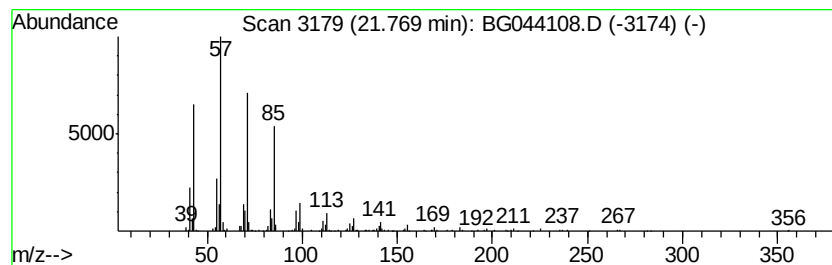
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ClientSampleId :
QVD-GW-01-45

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 7 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.77	2.78 ng	184955	Chrysene-d12		21.56
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044108.D
Acq On : 20 Jan 2020 21:38
Operator : JU
Sample : L1165-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

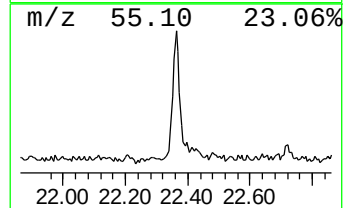
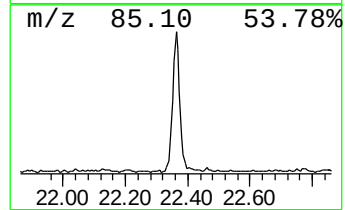
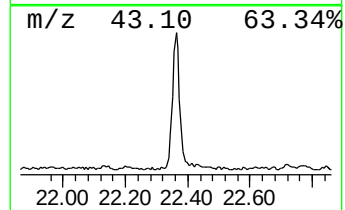
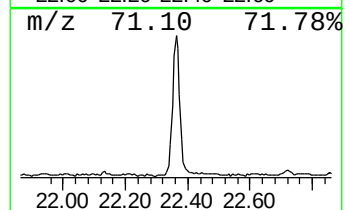
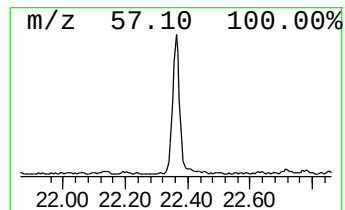
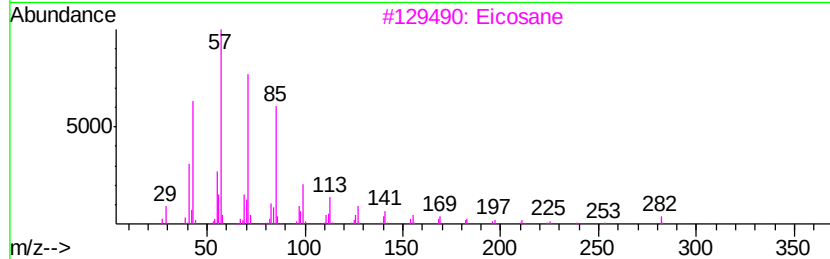
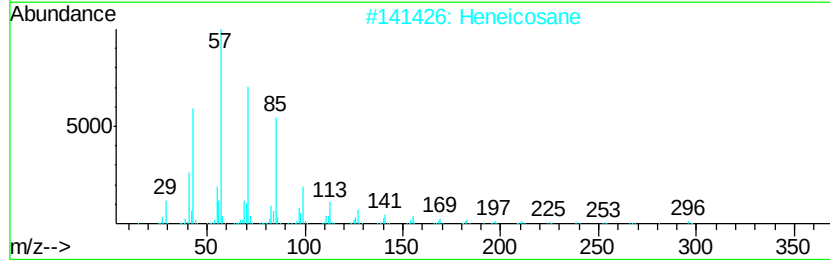
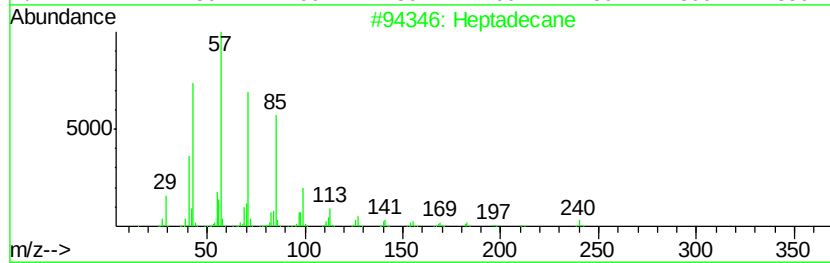
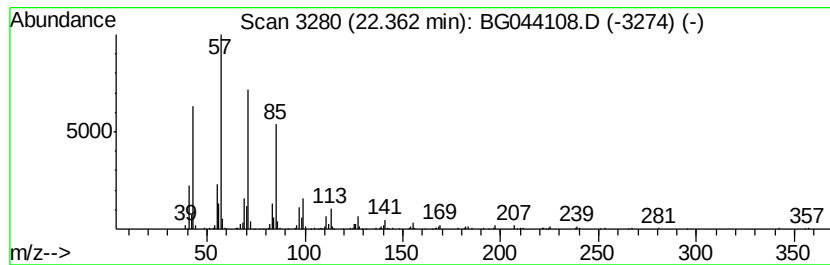
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ClientSampleId :
QVD-GW-01-45

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 8 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.36	4.79 ng	318889	Chrysene-d12	21.56	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Heneicosane	296	C21H44	000629-94-7	91
3	Eicosane	282	C20H42	000112-95-8	91
4	Hexacosane	366	C26H54	000630-01-3	91
5	Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044108.D
Acq On : 20 Jan 2020 21:38
Operator : JU
Sample : L1165-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

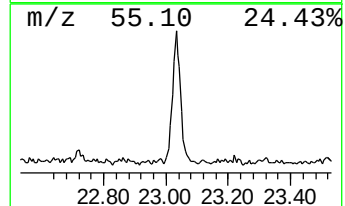
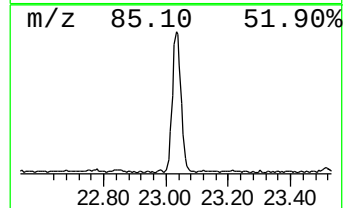
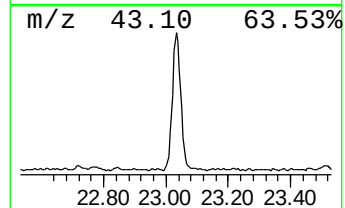
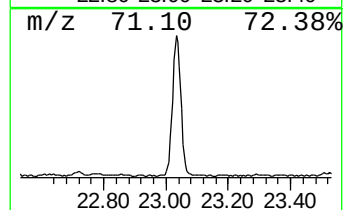
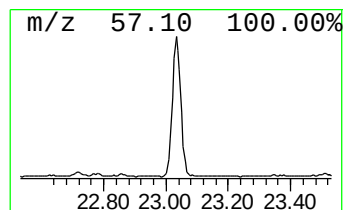
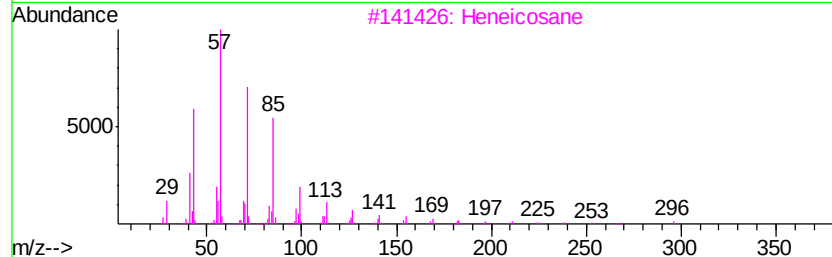
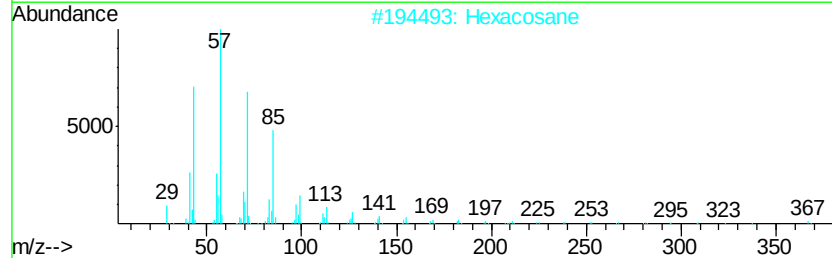
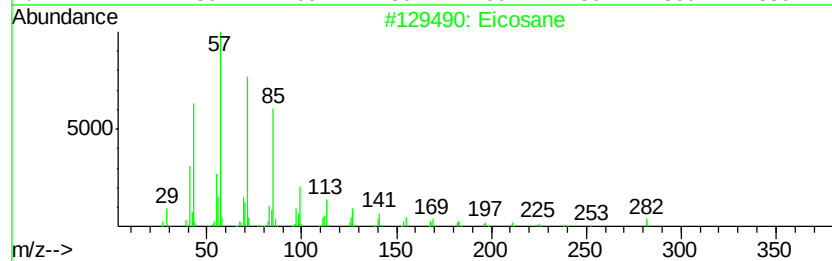
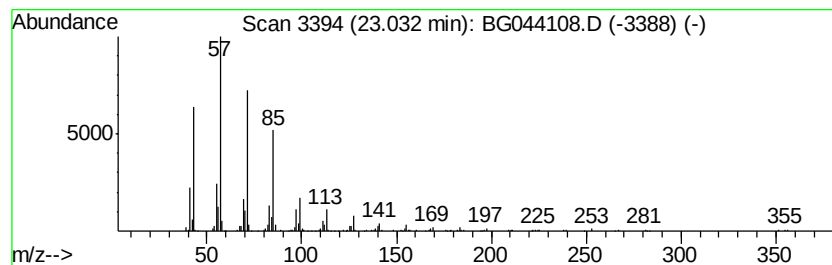
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane, 9-octyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
23.03	6.57 ng	437436	Chrysene-d12		21.56
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	98
2	Hexacosane	366	C26H54	000630-01-3	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
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Acq On : 20 Jan 2020 21:38
Operator : JU
Sample : L1165-01
Misc :
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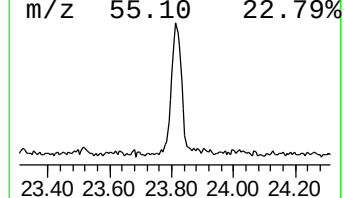
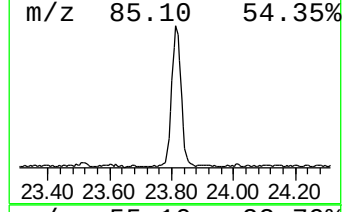
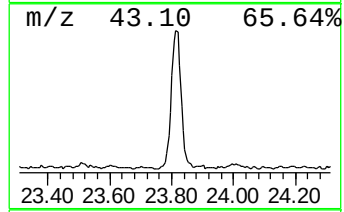
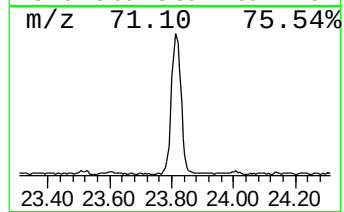
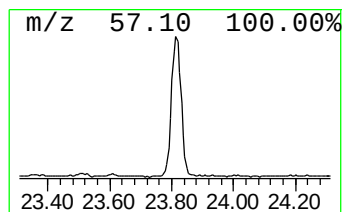
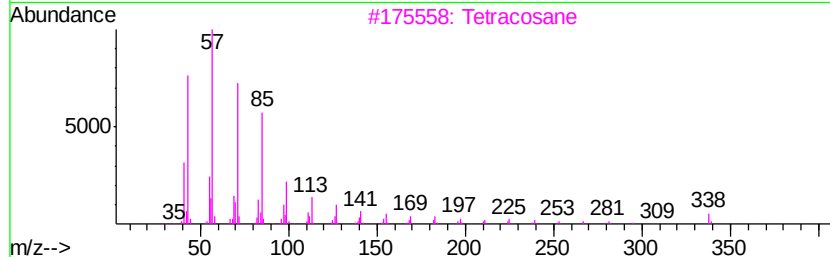
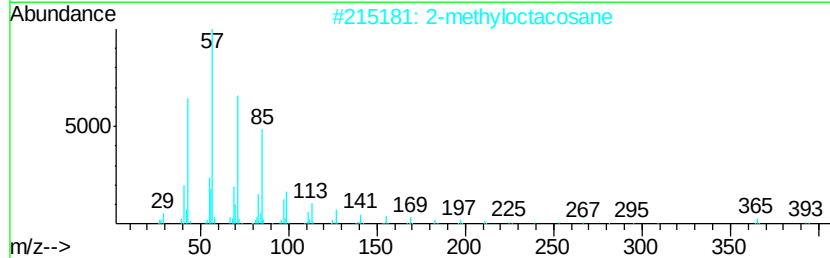
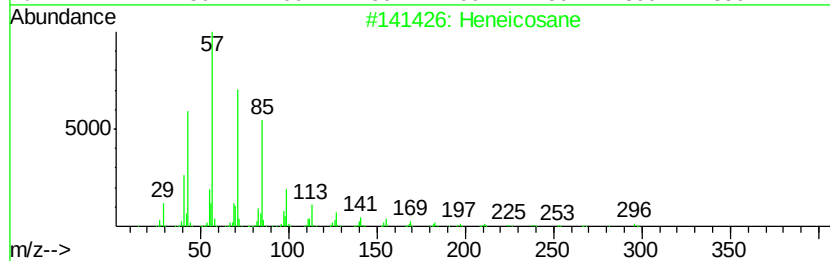
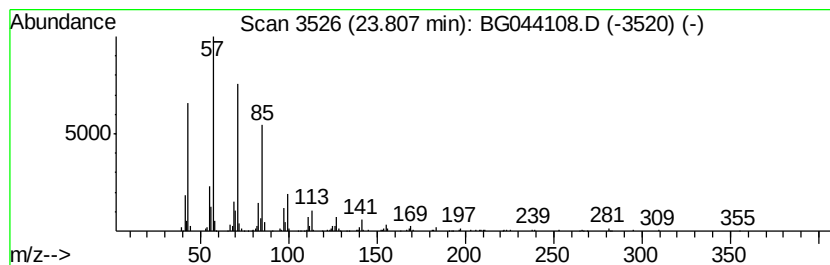
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.81	6.87 ng	475664	Perylene-d12	24.67	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	2-methyloctacosane	408	C29H60	1000376-72-8	91
3	Tetracosane	338	C24H50	000646-31-1	90
4	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5	Tetradecane	198	C14H30	000629-59-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044108.D
Acq On : 20 Jan 2020 21:38
Operator : JU
Sample : L1165-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

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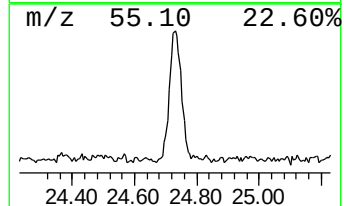
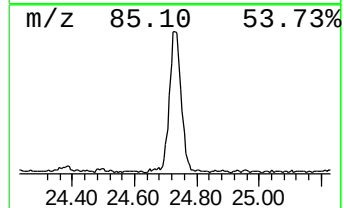
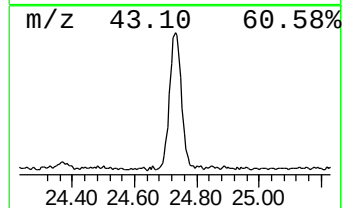
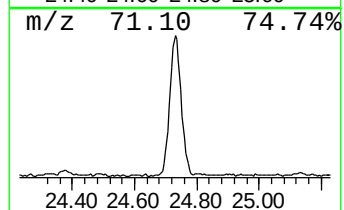
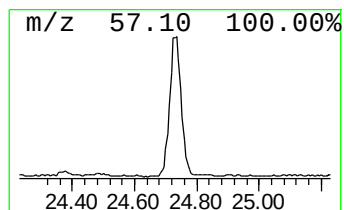
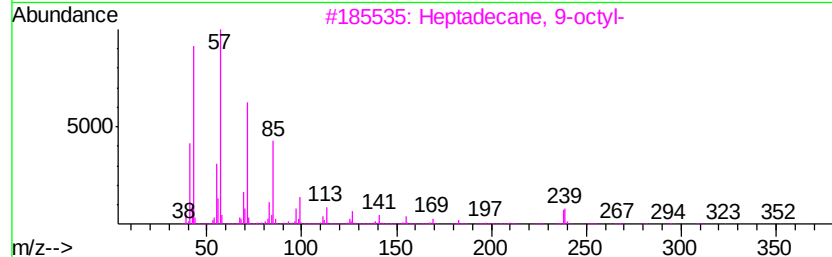
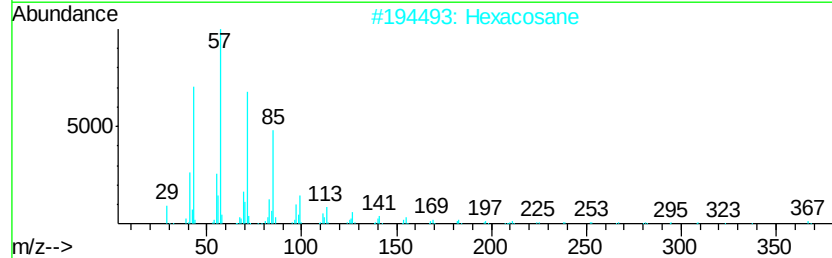
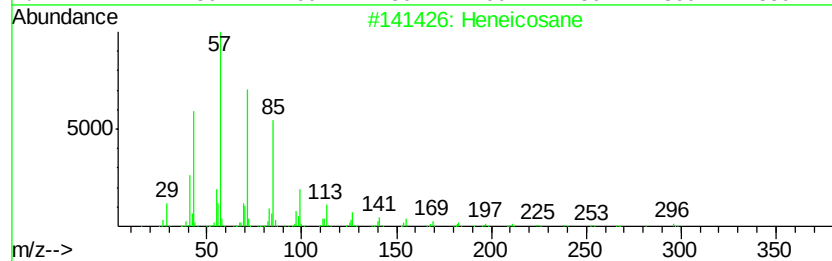
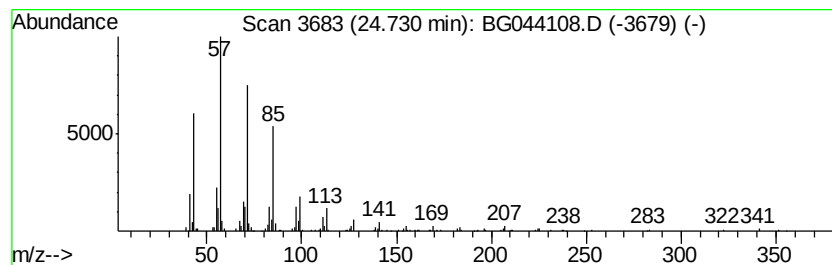
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.73	6.69 ng	463406	Perylene-d12	24.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044108.D
Acq On : 20 Jan 2020 21:38
Operator : JU
Sample : L1165-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
QVD-GW-01-45

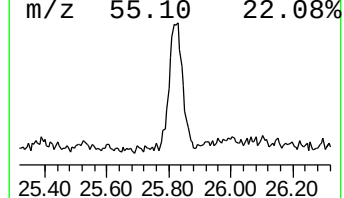
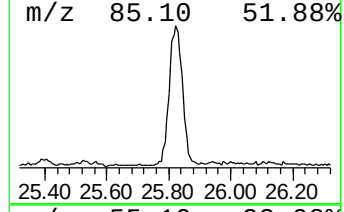
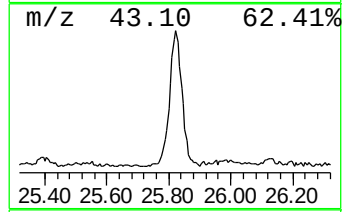
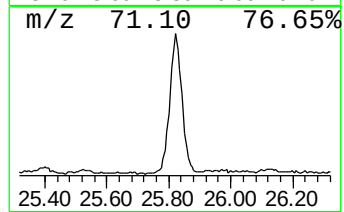
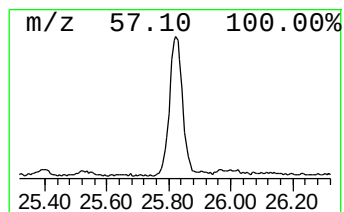
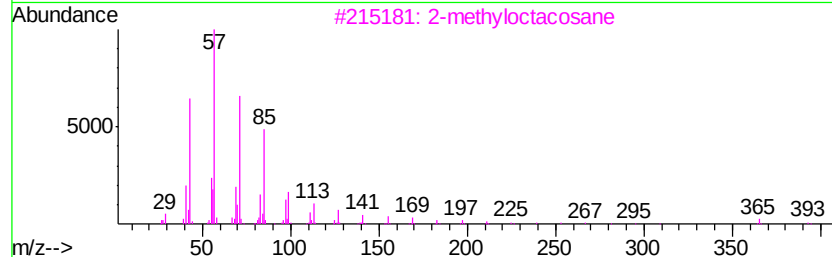
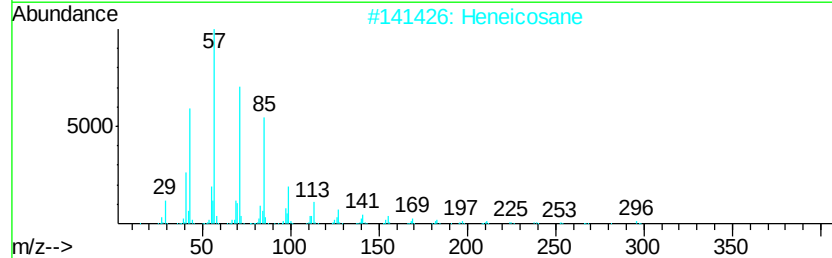
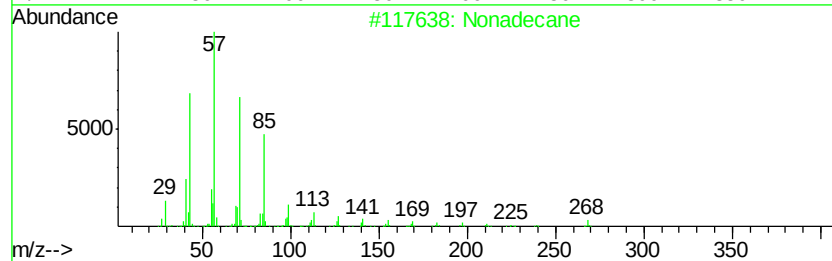
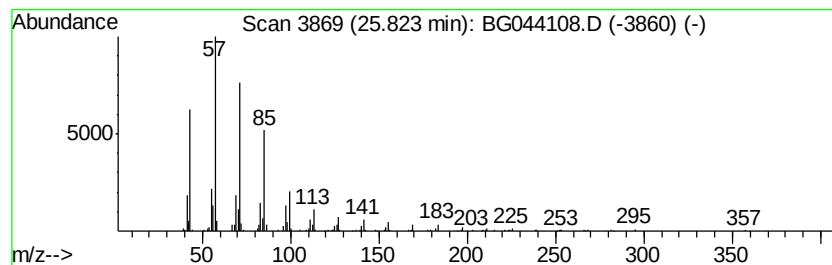
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.82	5.71 ng	395069	Perylene-d12	24.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044108.D
Acq On : 20 Jan 2020 21:38
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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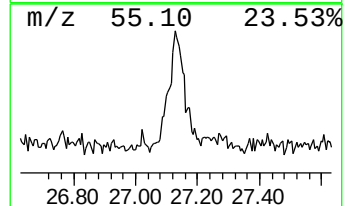
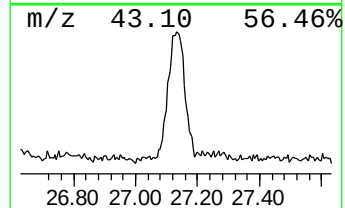
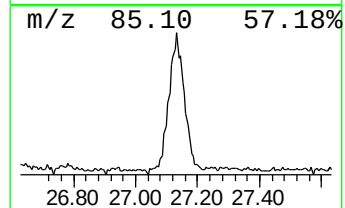
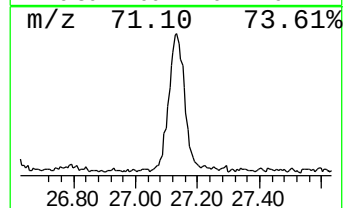
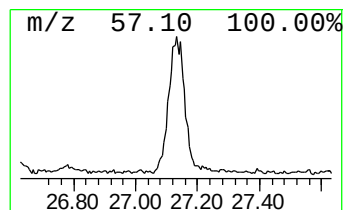
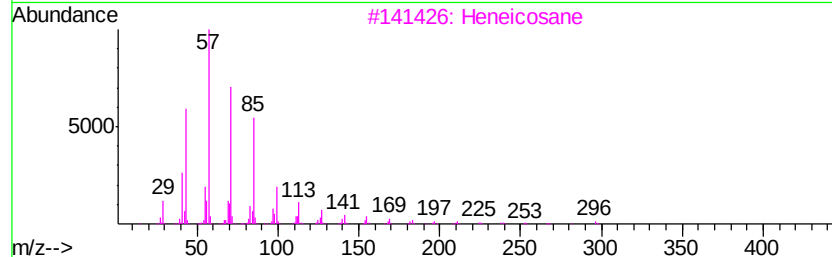
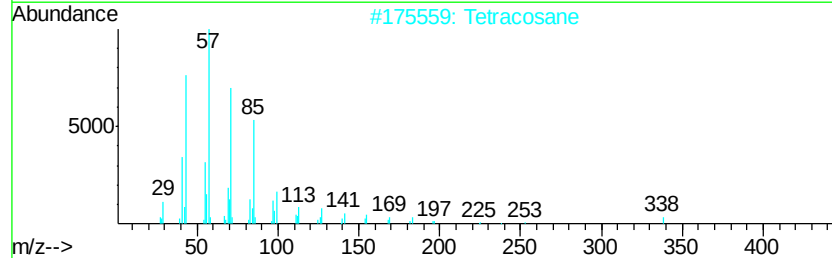
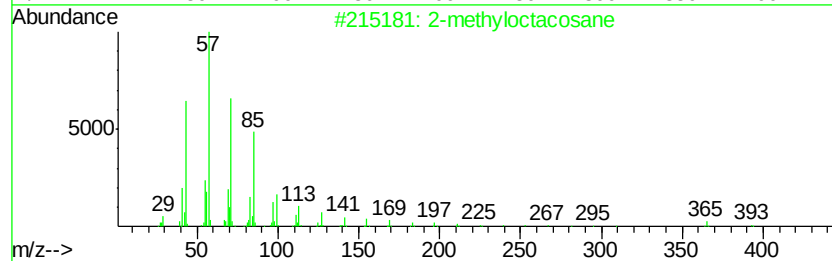
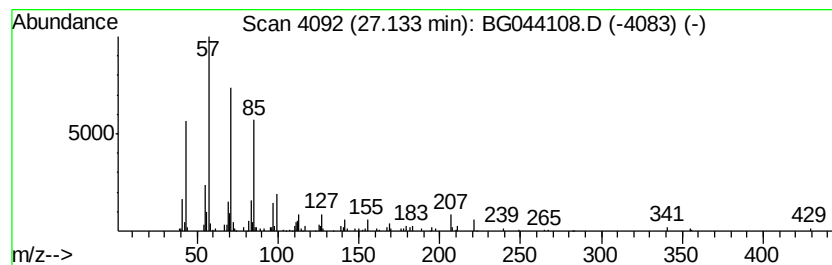
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-methyloctacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.13	2.98 ng	206233	Perylene-d12	24.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	87
2		Tetracosane	338	C24H50	000646-31-1	87
3		Heneicosane	296	C21H44	000629-94-7	86
4		Octacosane	394	C28H58	000630-02-4	86
5		Nonadecane	268	C19H40	000629-92-5	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG012020\
 Data File : BG044108.D
 Acq On : 20 Jan 2020 21:38
 Operator : JU
 Sample : L1165-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 QVD-GW-01-45

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
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Ethanol, 2-butoxy-	6.01	5.8	ng	130828	1	7.93	448308	20.0
unknown7.47	7.47	89.7	ng	2009880	1	7.93	448308	20.0
2,4,7,9-Tetrameth...	13.47	2.2	ng	116795	3	14.57	1070180	20.0
Oxalic acid, isob...	21.23	7.5	ng	502377	5	21.56	1331810	20.0
Eicosane	21.77	2.8	ng	184955	5	21.56	1331810	20.0
Heptadecane	22.36	4.8	ng	318889	5	21.56	1331810	20.0
Heptadecane, 9-oc...	23.03	6.6	ng	437436	5	21.56	1331810	20.0
Tetracosane	23.81	6.9	ng	475664	6	24.67	1384470	20.0
Pentacosane	24.73	6.7	ng	463406	6	24.67	1384470	20.0
Nonadecane	25.82	5.7	ng	395069	6	24.67	1384470	20.0
2-methyloctacosane	27.13	3.0	ng	206233	6	24.67	1384470	20.0