

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
 Data File : BG055651.D
 Acq On : 16 Nov 2022 17:38
 Operator : CG/JU
 Sample : N5607-13
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-40B

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055651.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.304	337	343	350	rBV	106735	164908	5.74%	0.995%
2	5.780	417	424	441	rVB	752973	1212017	42.22%	7.313%
3	7.390	688	698	712	rBV	784426	1347301	46.93%	8.130%
4	8.253	838	845	852	rVB	136950	228454	7.96%	1.378%
5	9.422	1036	1044	1053	rVB	460911	809342	28.19%	4.884%
6	10.833	1276	1284	1297	rBV2	48057	113706	3.96%	0.686%
7	11.079	1319	1326	1334	rBV	204205	356168	12.41%	2.149%
8	13.500	1730	1738	1747	rVB	1304954	2028831	70.67%	12.242%
9	14.187	1847	1855	1861	rBV2	30640	53310	1.86%	0.322%
10	14.475	1896	1904	1909	rBV10	25112	66255	2.31%	0.400%
11	14.875	1962	1972	1979	rVB	338562	539123	18.78%	3.253%
12	15.180	2018	2024	2030	rVB	41298	65593	2.28%	0.396%
13	15.797	2124	2129	2134	rBV3	20409	30882	1.08%	0.186%
14	16.355	2217	2224	2230	rBV	1087206	1668313	58.11%	10.067%
15	16.567	2254	2260	2268	rVB4	32540	61170	2.13%	0.369%
16	17.061	2334	2344	2349	rBV10	29844	99595	3.47%	0.601%
17	17.119	2349	2354	2360	rVB	131176	193500	6.74%	1.168%
18	17.613	2432	2438	2444	rBV	431392	646014	22.50%	3.898%
19	18.477	2576	2585	2602	rBV	307511	553291	19.27%	3.339%
20	19.734	2795	2799	2811	rBV	309626	450705	15.70%	2.720%
21	20.204	2873	2879	2885	rVB	2165581	2870743	100.00%	17.322%
22	20.486	2924	2927	2933	rBV	32785	46133	1.61%	0.278%
23	20.756	2969	2973	2978	rBV	32493	48900	1.70%	0.295%
24	20.991	3009	3013	3016	rBV2	42924	49913	1.74%	0.301%
25	21.438	3085	3089	3093	rBV	37966	58872	2.05%	0.355%
26	21.520	3099	3103	3114	rBV3	59211	137015	4.77%	0.827%
27	21.908	3163	3169	3181	rVB2	428344	749865	26.12%	4.525%
28	22.090	3196	3200	3207	rVB	57736	88169	3.07%	0.532%
29	22.202	3215	3219	3227	rBV3	43213	85381	2.97%	0.515%
30	22.730	3304	3309	3316	rVB	75470	127933	4.46%	0.772%
31	23.465	3428	3434	3444	rVB	68127	143648	5.00%	0.867%
32	24.329	3574	3581	3590	rVB	87943	214176	7.46%	1.292%
33	24.546	3613	3618	3625	rVB7	34843	79138	2.76%	0.478%
34	25.327	3740	3751	3769	rVB2	280935	963114	33.55%	5.811%
35	26.555	3953	3960	3973	rVB3	66456	221221	7.71%	1.335%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

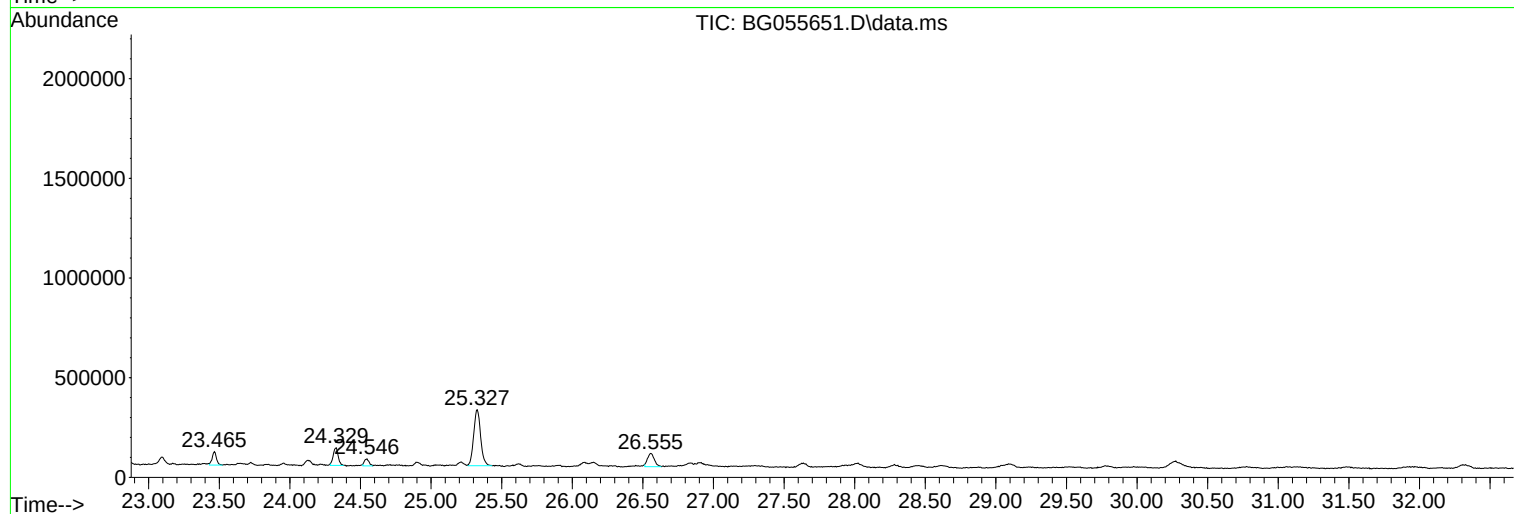
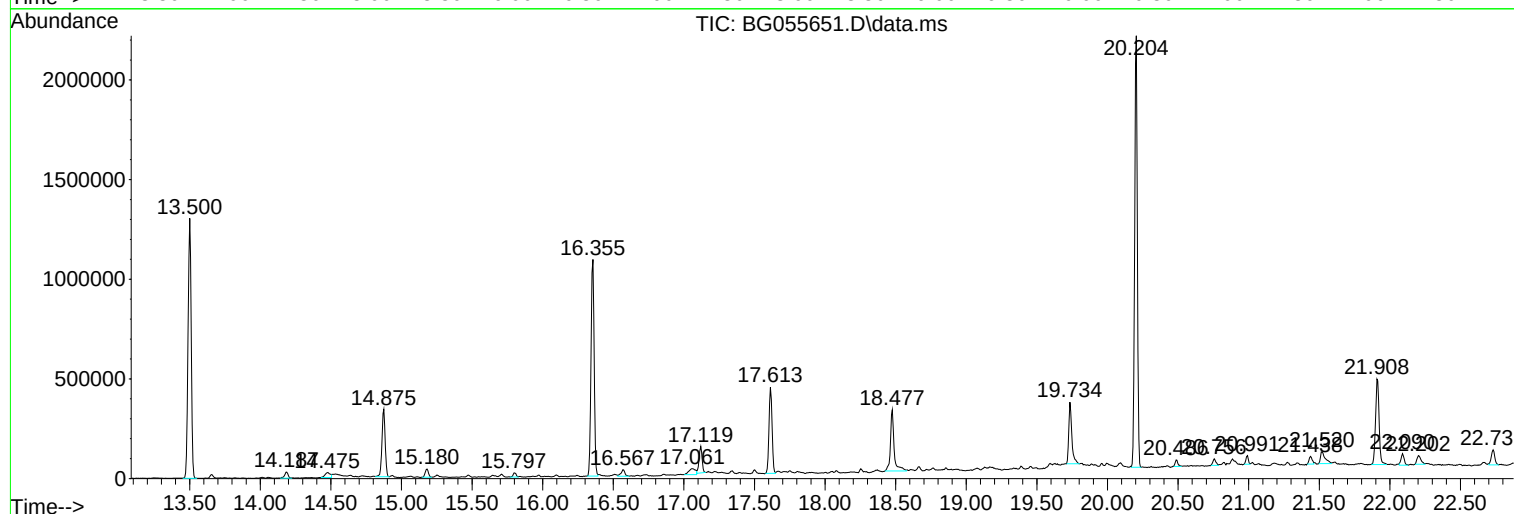
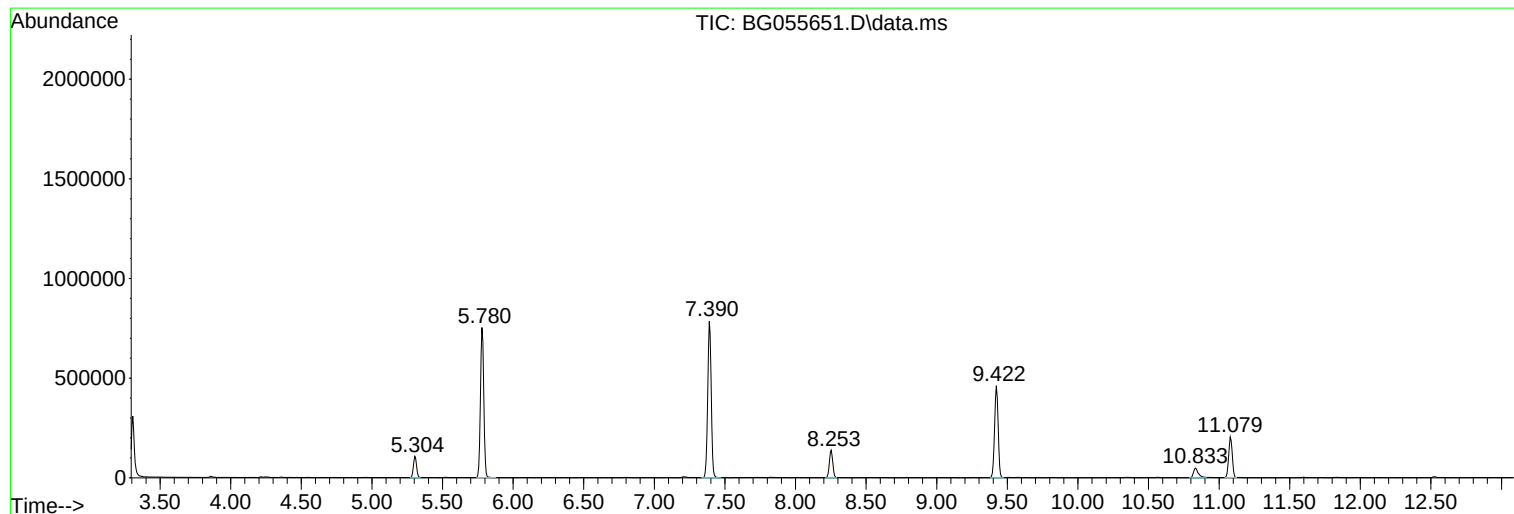
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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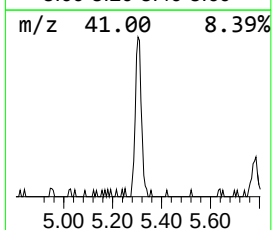
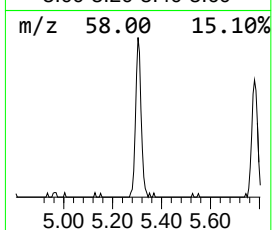
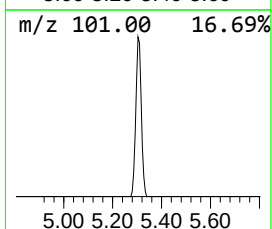
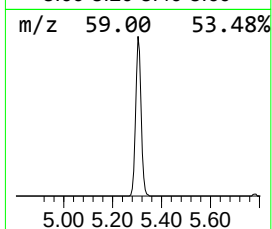
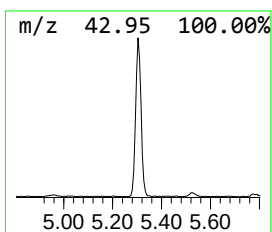
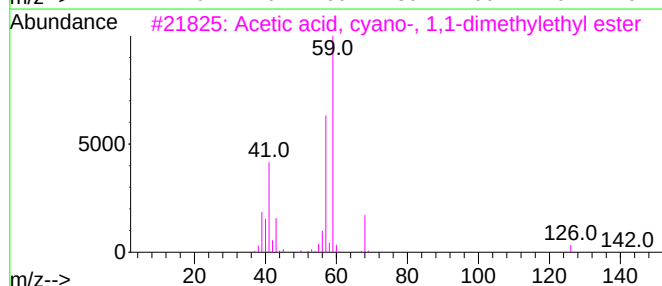
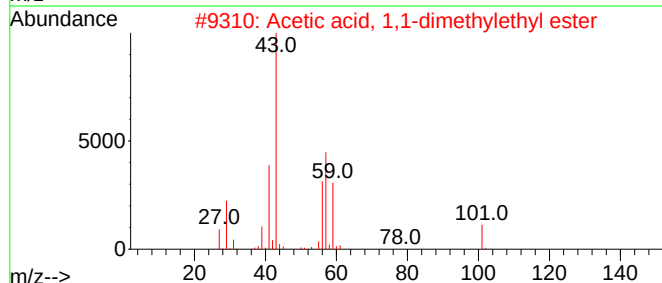
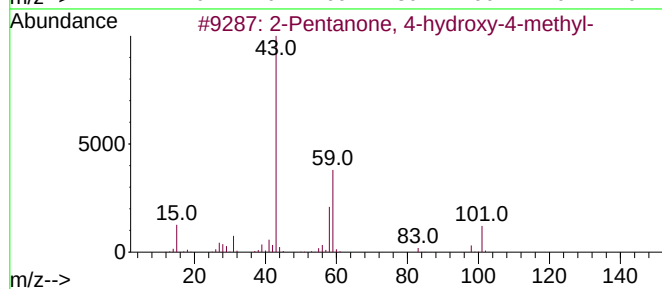
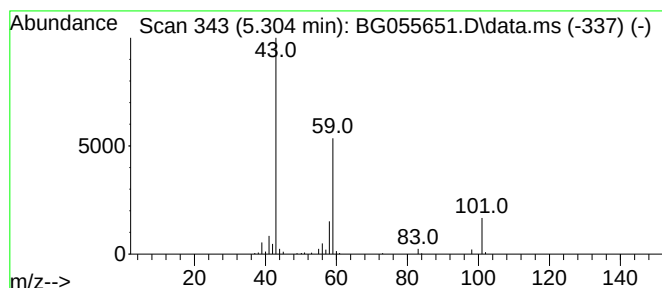
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.304	14.44 ng	164908	1,4-Dichlorobenzene-d4	8.253

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		
5	Butane	58	C4H10	000106-97-8	9		



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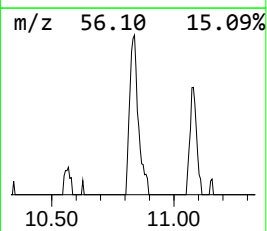
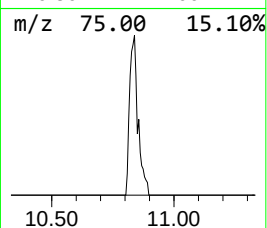
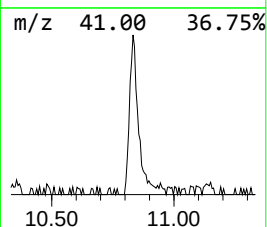
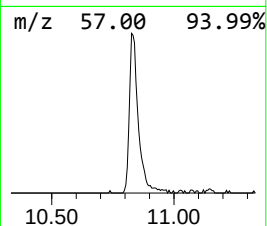
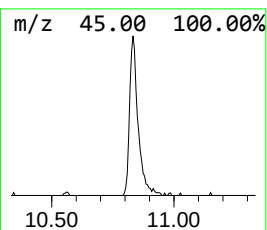
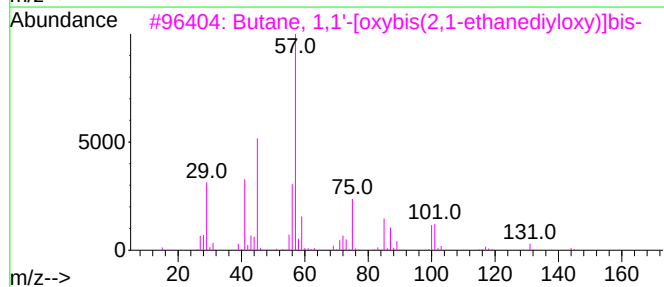
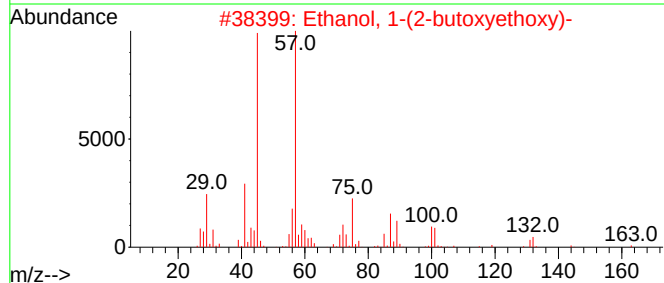
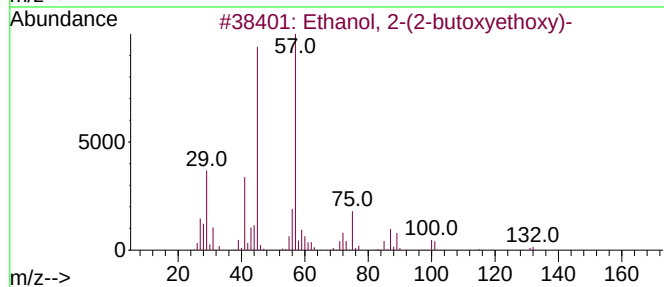
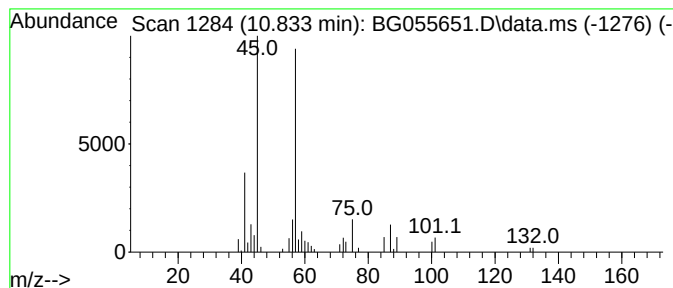
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.833	6.38 ng	113706	Naphthalene-d8	11.079

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	64
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
5			Di-sec-Butyl ether	130	C8H18O	006863-58-7	53



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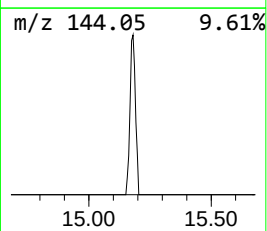
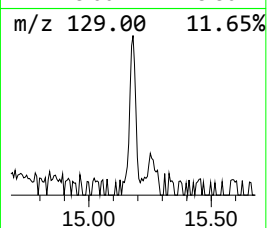
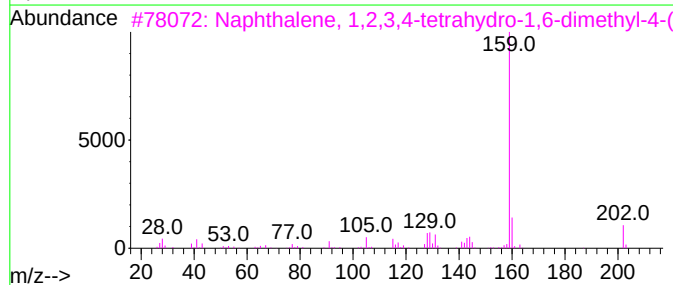
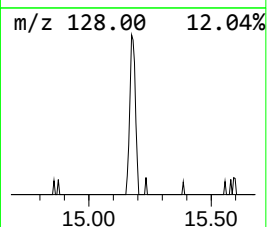
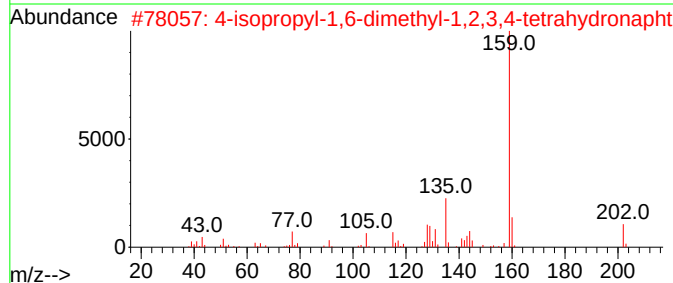
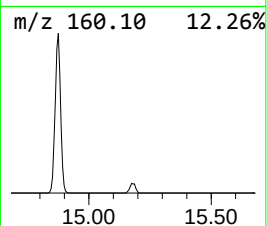
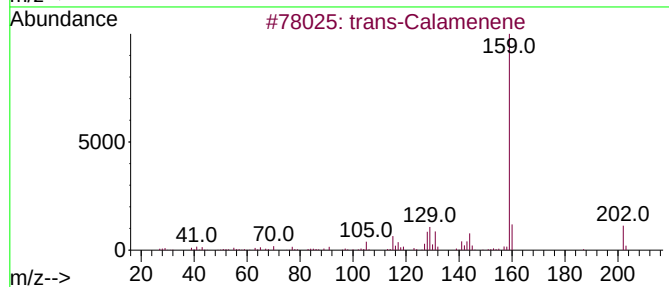
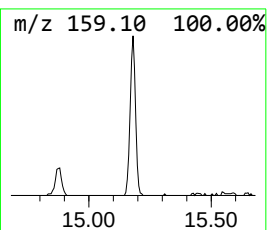
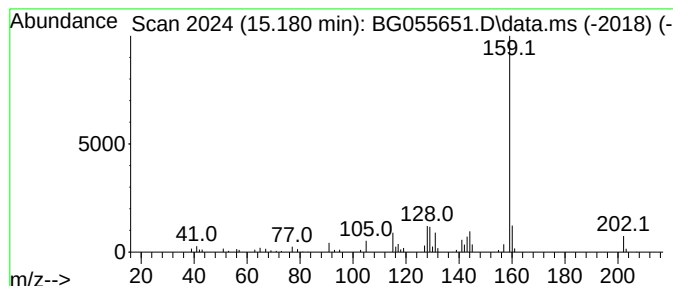
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 trans-Calamenene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.180	2.43 ng	65593	Acenaphthene-d10	14.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Calamenene	202	C15H22	073209-42-4	97
2			4-isopropyl-1,6-dimethyl-1,2,3,4...	202	C15H22	1000378-99-6	97
3			Naphthalene, 1,2,3,4-tetrahydro...	202	C15H22	000483-77-2	95
4			cis-Calamenene	202	C15H22	072937-55-4	94
5			Naphthalene, 1,2,3,4-tetrahydro...	174	C13H18	000475-03-6	86



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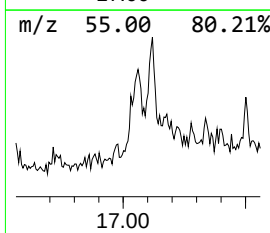
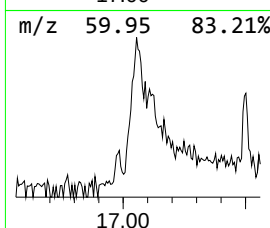
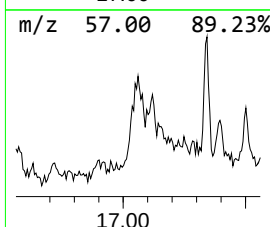
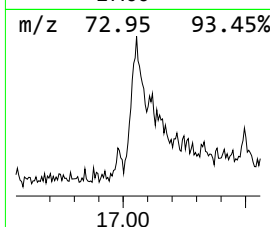
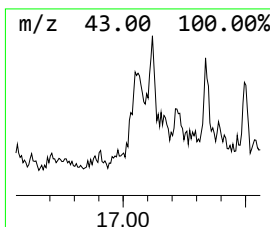
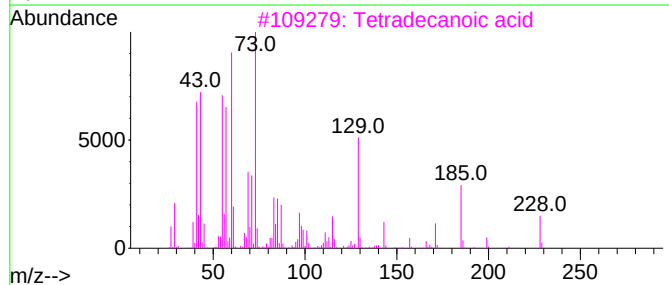
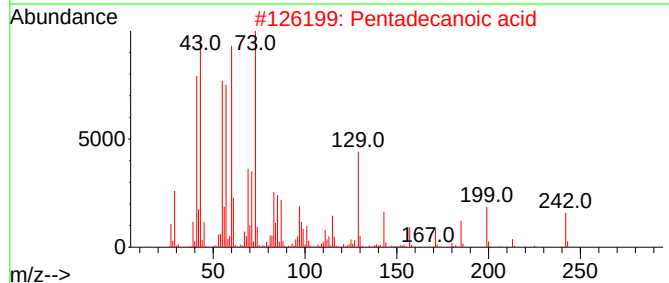
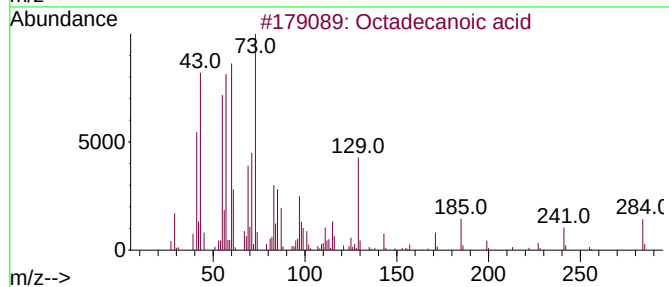
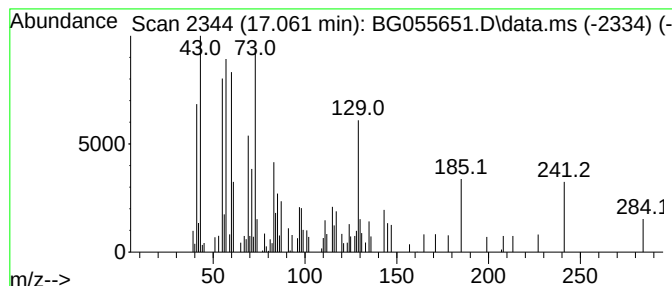
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.061	3.08 ng	99595	Phenanthrene-d10	17.613

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	96
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	52
5			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	46



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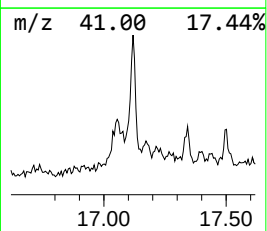
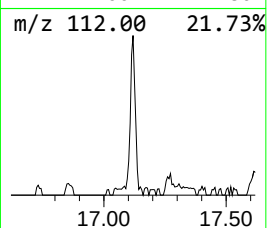
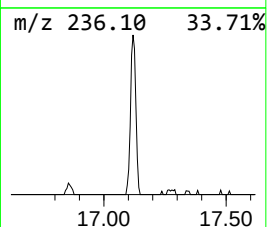
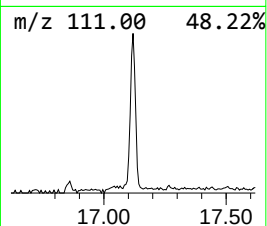
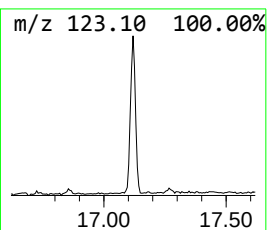
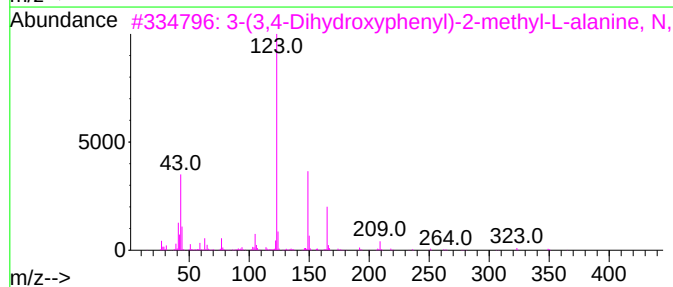
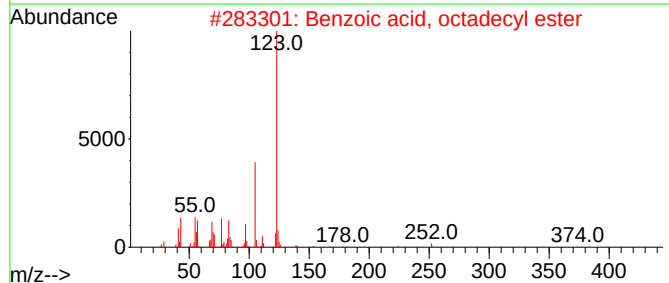
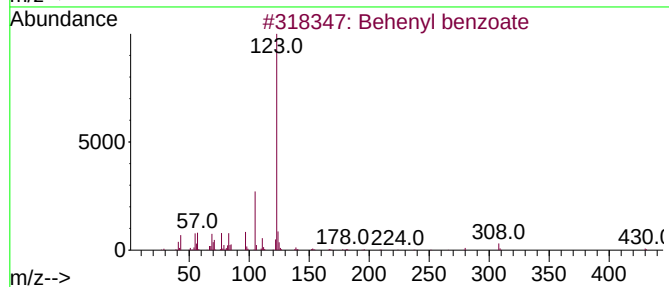
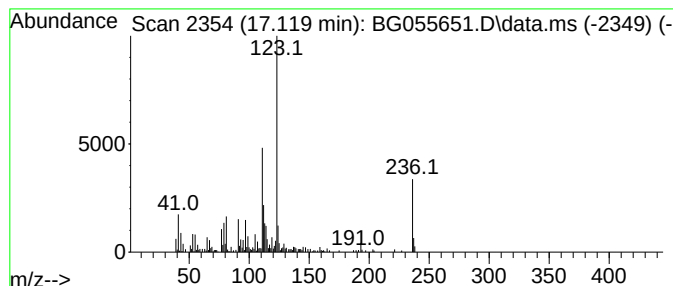
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Peak Number 6 unknown17.119 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.119	5.99 ng	193500	Phenanthrene-d10	17.613

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenyl benzoate	430	C29H50O2	103403-38-9	38
2			Benzoic acid, octadecyl ester	374	C25H42O2	010578-34-4	38
3			3-(3,4-Dihydroxyphenyl)-2-methyl...	483	C23H33NO10	1000453-24-7	35
4			5-Amino-1,3,3-trimethylcyclohexa...	362	C14H20F6N2O2	1000454-06-2	35
5			Cyclopent-2-ene-1-carboxylic aci...	196	C12H20O2	1000195-65-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
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Acq On : 16 Nov 2022 17:38
Operator : CG/JU
Sample : N5607-13
Misc :
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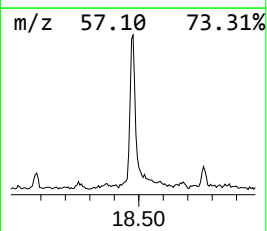
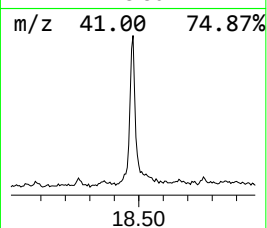
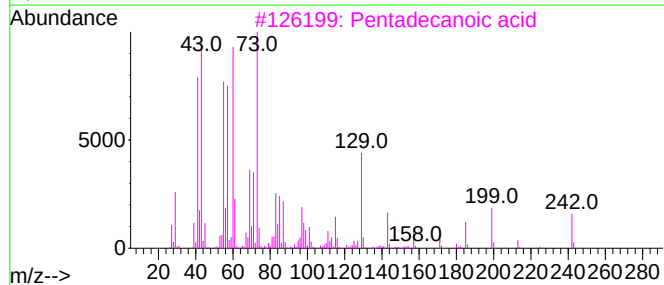
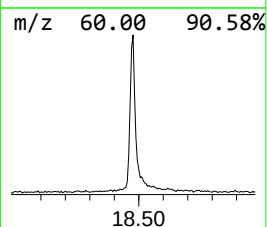
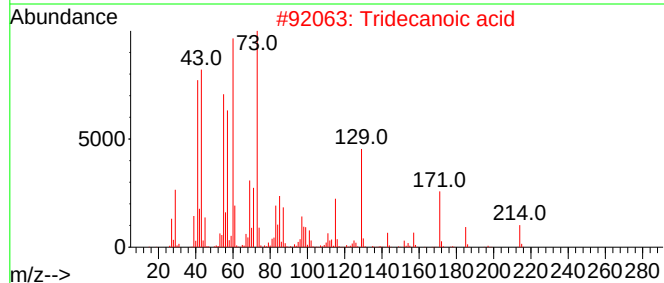
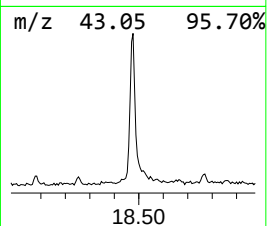
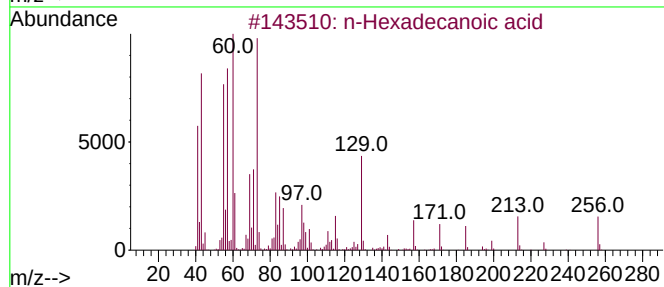
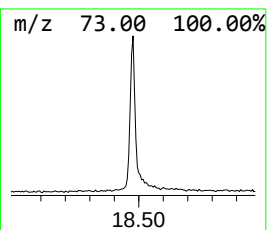
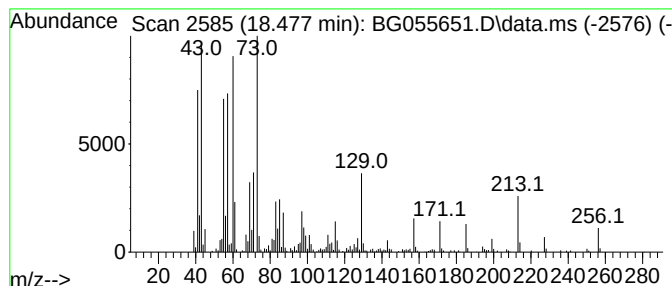
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.477	17.13 ng	553291	Phenanthrene-d10		17.613	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	99
2	Tridecanoic acid		214	C13H26O2	000638-53-9	93
3	Pentadecanoic acid		242	C15H30O2	001002-84-2	90
4	Tetradecanoic acid		228	C14H28O2	000544-63-8	72
5	Undecanoic acid		186	C11H22O2	000112-37-8	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
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Sample : N5607-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

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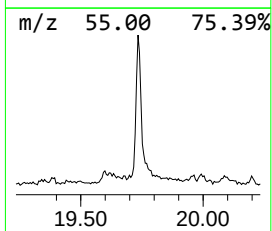
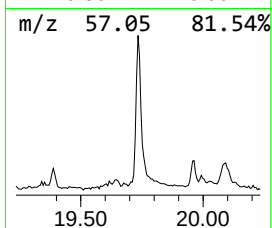
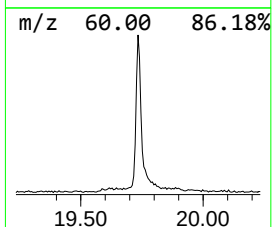
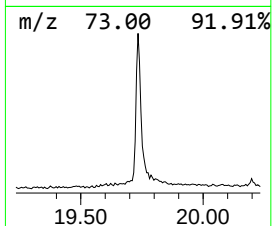
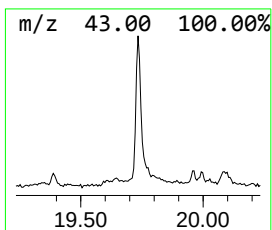
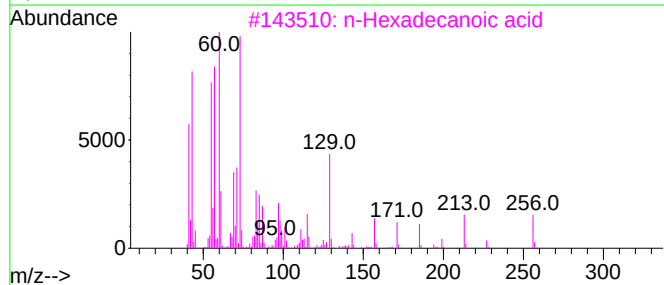
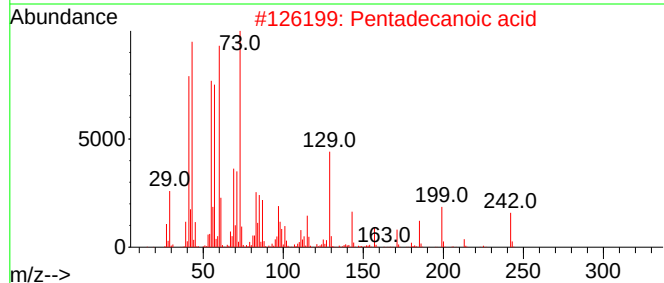
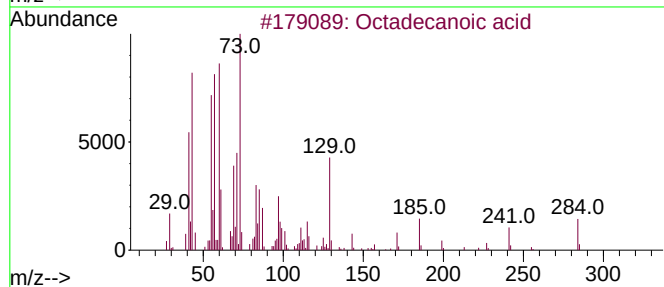
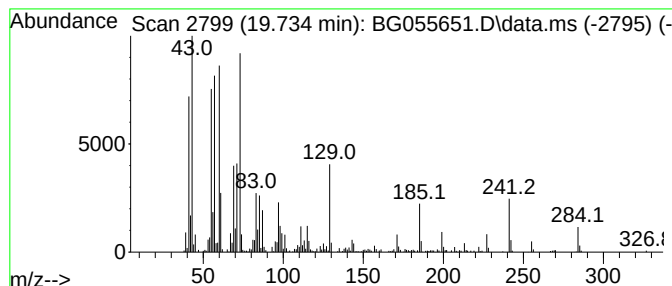
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pentadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.734	13.95 ng	450705	Phenanthrene-d10	17.613

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	95
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
Operator : CG/JU
Sample : N5607-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-40B

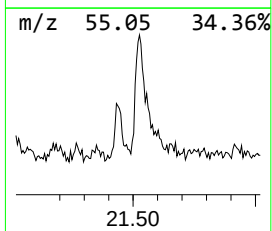
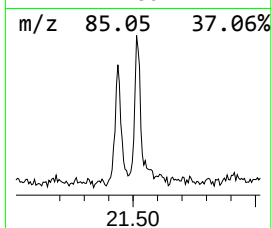
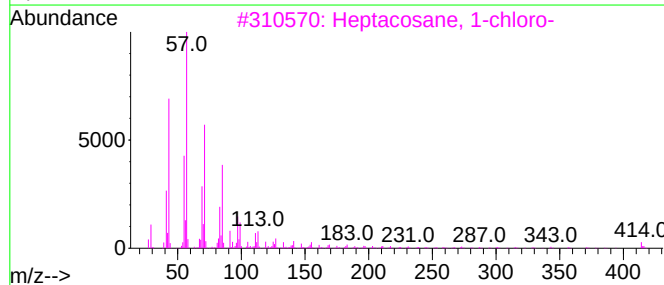
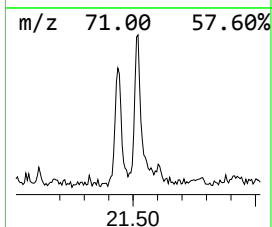
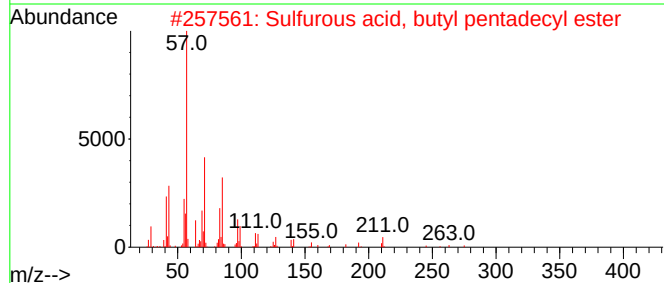
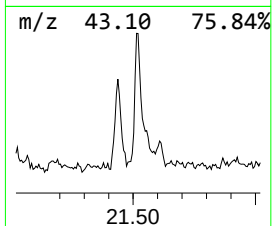
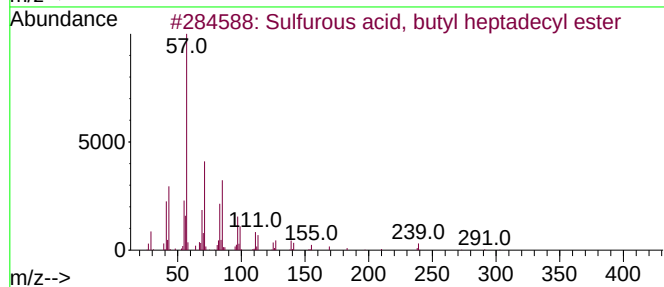
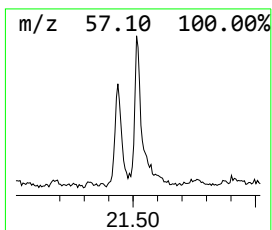
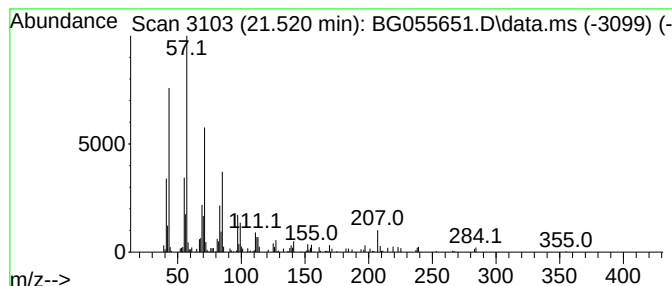
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Sulfurous acid, butyl hepta... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.520	3.65 ng	137015	Chrysene-d12	21.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	90
2			Sulfurous acid, butyl pentadecyl...	348	C19H40O3S	1000309-18-2	87
3			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	87
4			Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	87
5			1-Chloroeicosane	316	C20H41Cl	042217-02-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
Operator : CG/JU
Sample : N5607-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

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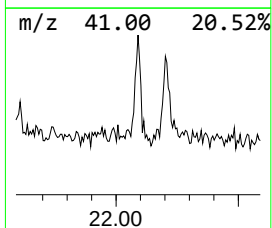
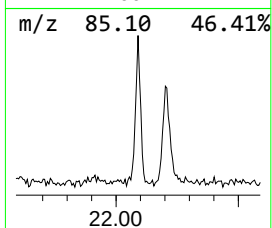
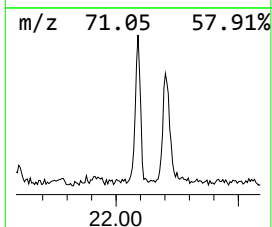
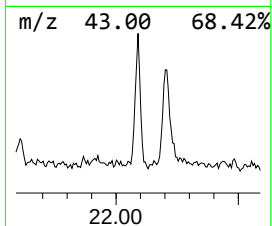
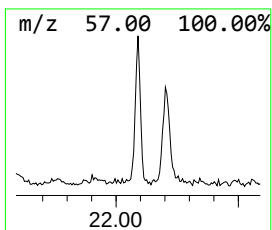
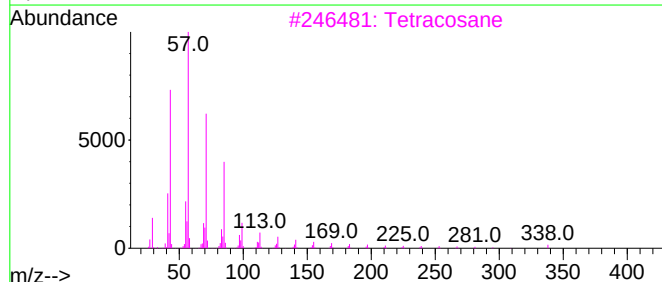
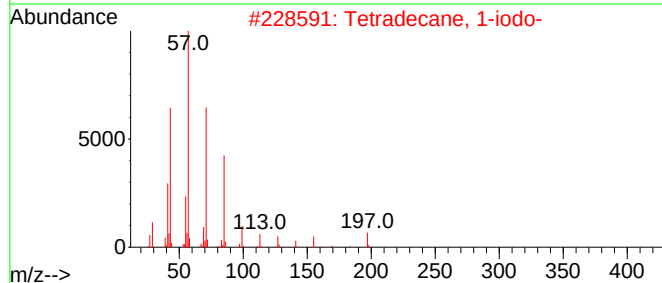
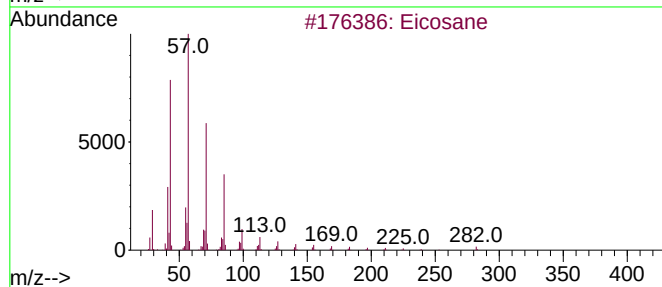
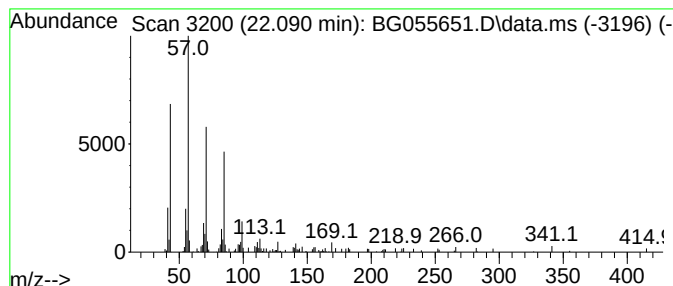
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.090	2.35 ng	88169	Chrysene-d12	21.908

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	96
2		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	83
3		Tetracosane	338	C24H50	000646-31-1	83
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
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Sample : N5607-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

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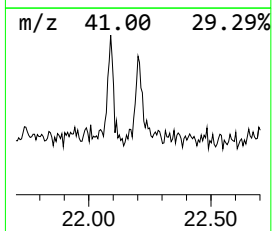
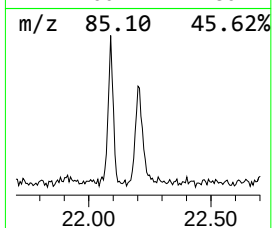
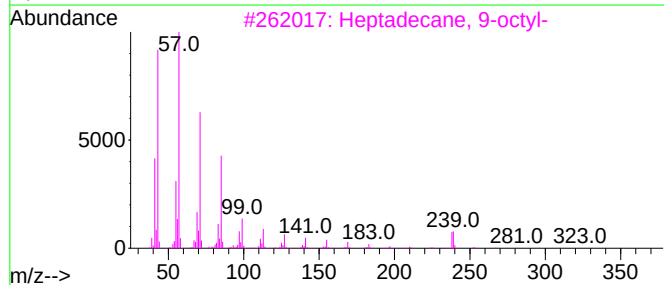
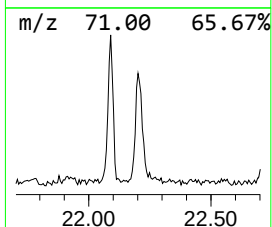
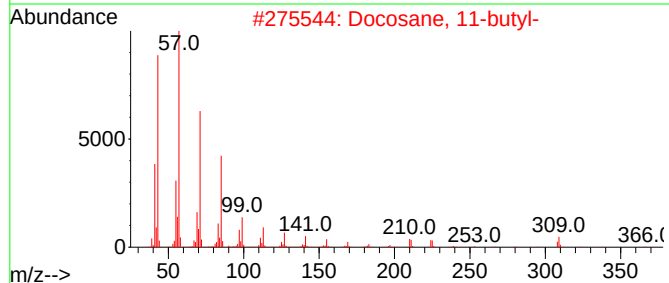
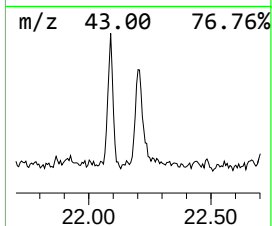
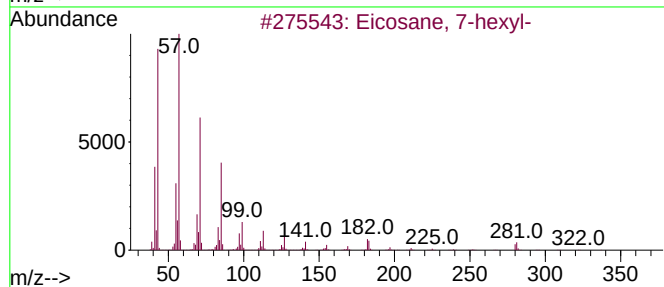
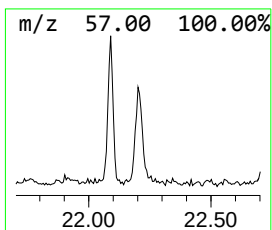
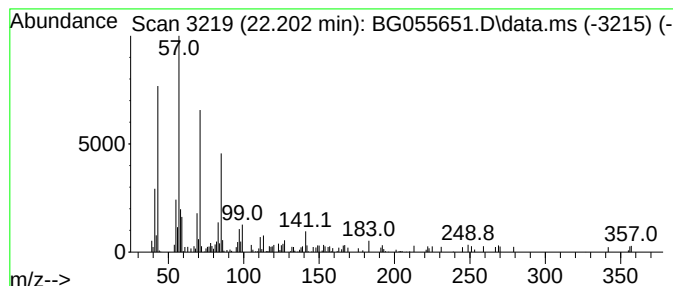
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Eicosane, 7-hexyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.202	2.28 ng	85381	Chrysene-d12	21.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	68
2			Docosane, 11-butyl-	366	C26H54	013475-76-8	68
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	64
4			4-Methylheneicosane	310	C22H46	025117-29-7	64
5			Hexatriacontane	507	C36H74	000630-06-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
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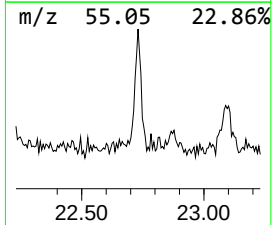
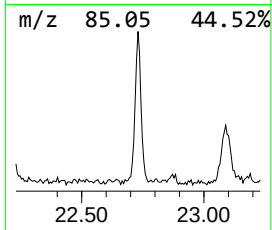
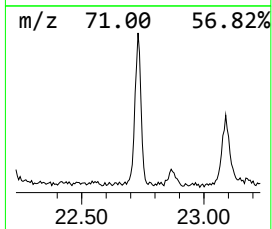
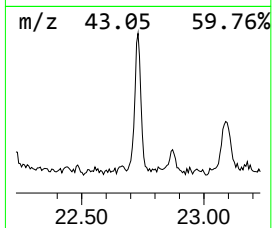
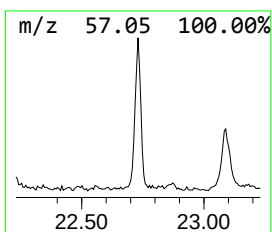
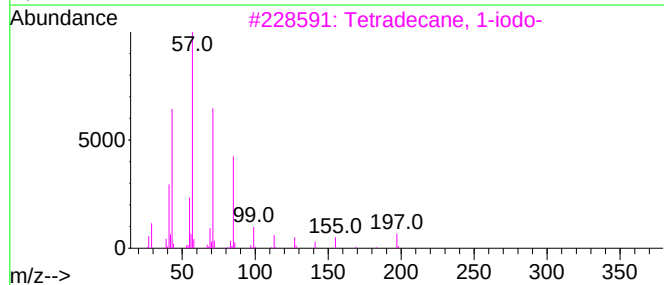
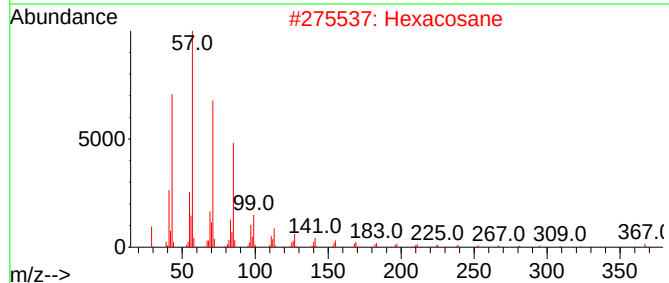
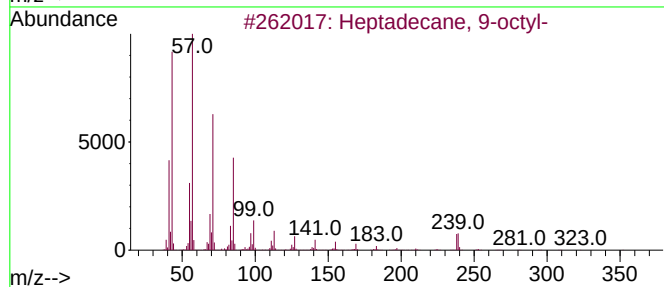
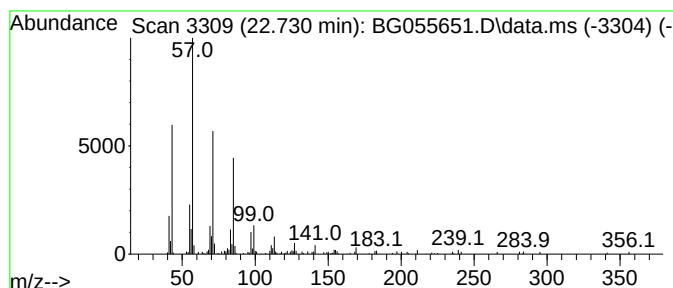
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecane, 9-octyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.730	3.41 ng	127933	Chrysene-d12	21.908	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
2	Hexacosane	366	C26H54	000630-01-3	87
3	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	83
4	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	83
5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
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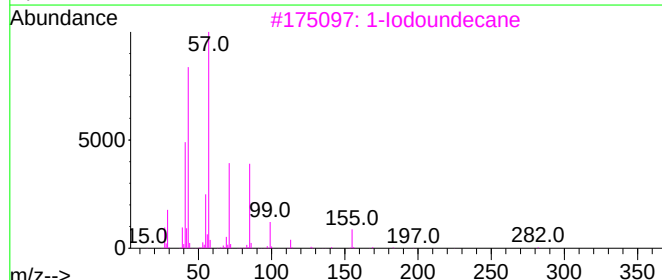
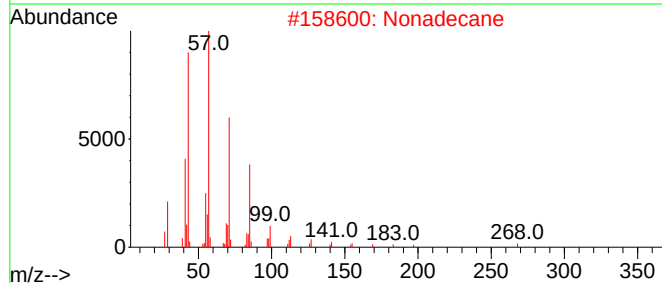
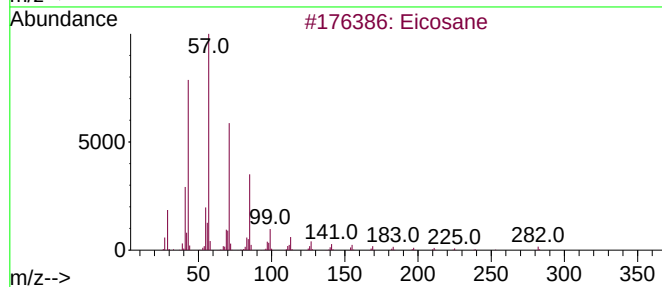
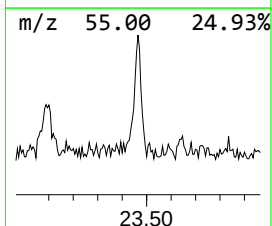
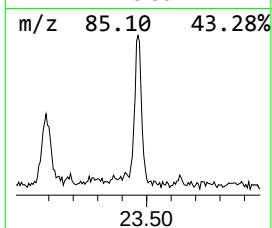
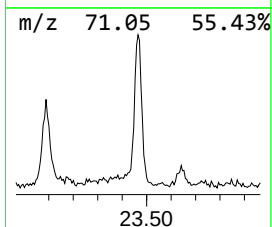
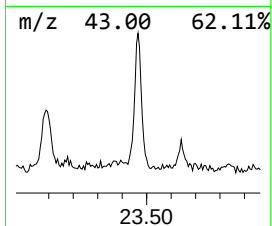
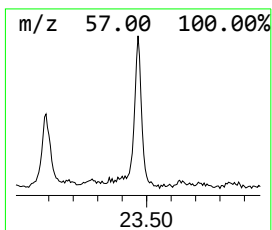
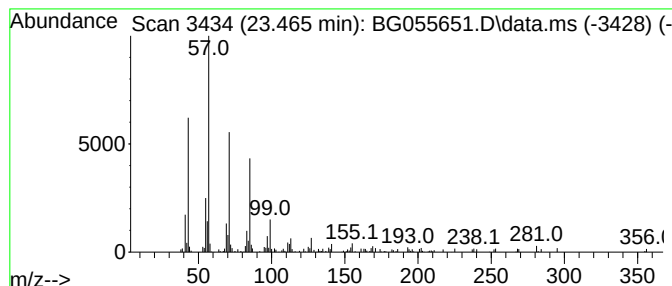
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.465	3.83 ng	143648	Chrysene-d12	21.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	95
2	Nonadecane			268	C19H40	000629-92-5	93
3	1-Iodoundecane			282	C11H23I	004282-44-4	87
4	Octacosane			394	C28H58	000630-02-4	83
5	Tridecane, 1-iodo-			310	C13H27I	035599-77-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
Operator : CG/JU
Sample : N5607-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-40B

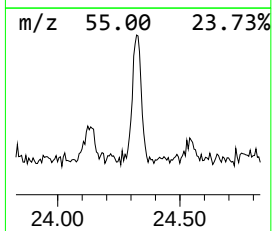
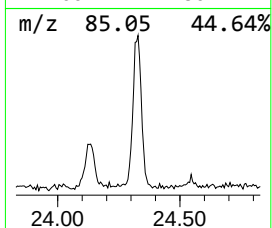
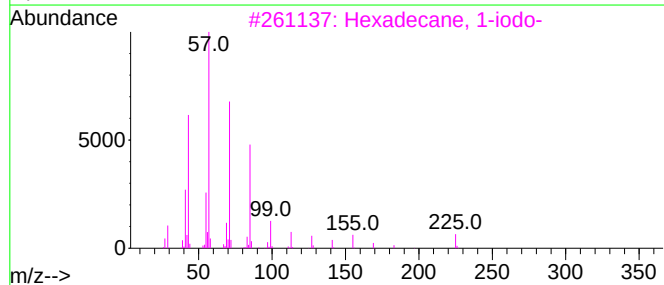
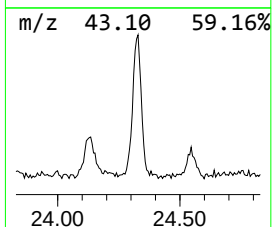
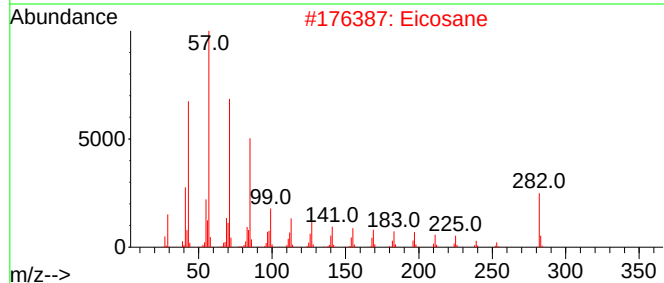
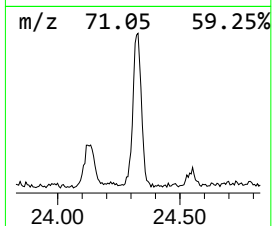
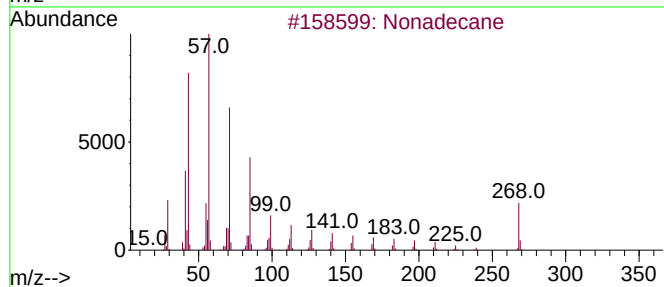
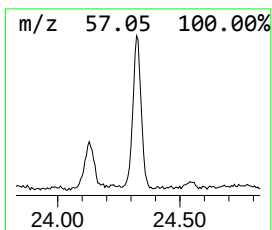
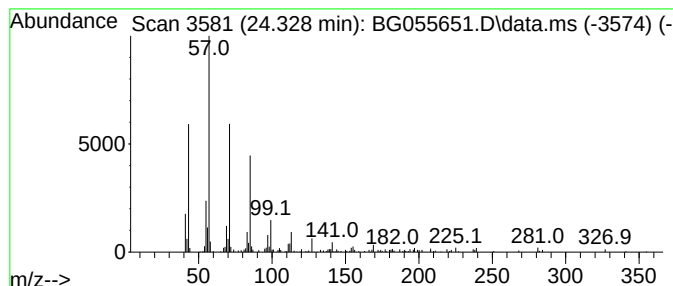
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Hexadecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.328	4.45 ng	214176	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	90
2			Eicosane	282	C20H42	000112-95-8	87
3			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
4			Heptacosane	380	C27H56	000593-49-7	87
5			Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
Operator : CG/JU
Sample : N5607-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

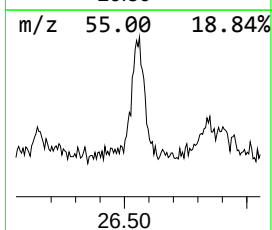
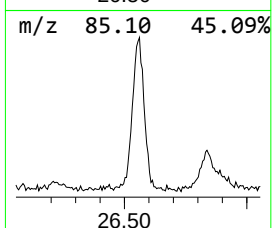
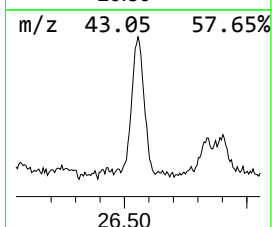
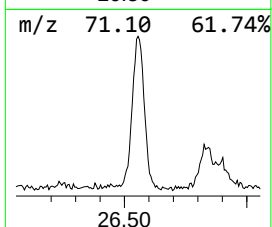
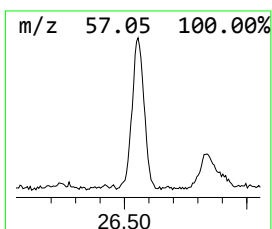
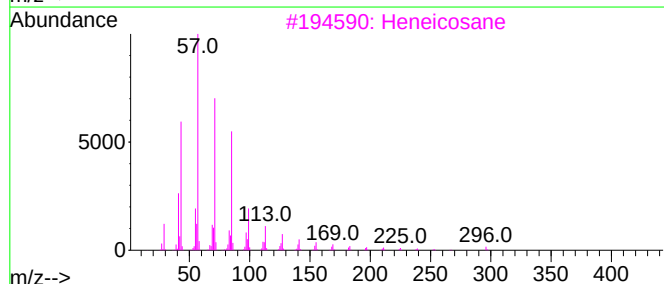
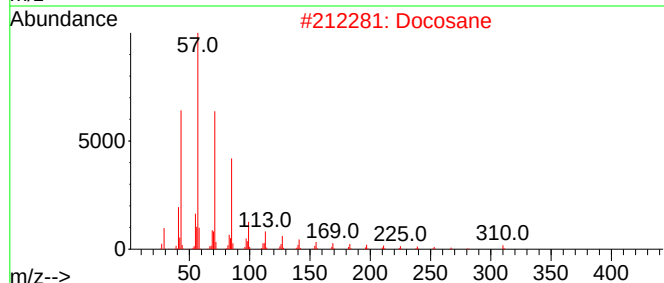
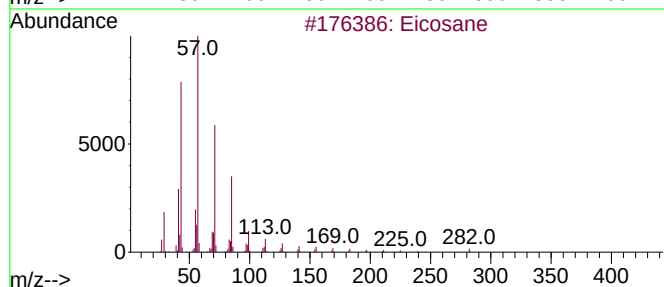
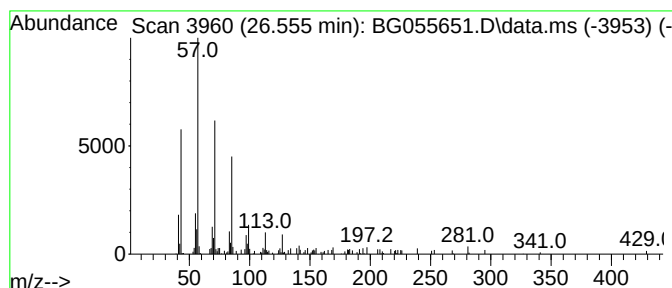
Instrument :
BNA_G
ClientSampleId :
SB-40B

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Docosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.555	4.59 ng	221221	Perylene-d12	25.321	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Docosane	310	C22H46	000629-97-0	87
3	Heneicosane	296	C21H44	000629-94-7	87
4	Nonadecane	268	C19H40	000629-92-5	86
5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
Data File : BG055651.D
Acq On : 16 Nov 2022 17:38
Operator : CG/JU
Sample : N5607-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-40B

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.304	14.4	ng	164908	1	8.253	228454	20.0
Ethanol, 2-(2-b...	10.833	6.4	ng	113706	2	11.079	356168	20.0
Naphthalene, de...	14.475	2.5	ng	66255	3	14.875	539123	20.0
trans-Calamenene	15.180	2.4	ng	65593	3	14.875	539123	20.0
Octadecanoic acid	17.061	3.1	ng	99595	4	17.613	646014	20.0
unknown17.119	17.119	6.0	ng	193500	4	17.613	646014	20.0
n-Hexadecanoic ...	18.477	17.1	ng	553291	4	17.613	646014	20.0
Pentadecanoic acid	19.734	13.9	ng	450705	4	17.613	646014	20.0
Sulfurous acid,...	21.520	3.6	ng	137015	5	21.908	749865	20.0
Eicosane	22.090	2.4	ng	88169	5	21.908	749865	20.0
Eicosane, 7-hexyl-	22.202	2.3	ng	85381	5	21.908	749865	20.0
Heptadecane, 9-...	22.730	3.4	ng	127933	5	21.908	749865	20.0
Nonadecane	23.465	3.8	ng	143648	5	21.908	749865	20.0
Hexadecane, 1-i...	24.328	4.5	ng	214176	6	25.321	963114	20.0
Docosane	26.555	4.6	ng	221221	6	25.321	963114	20.0