

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
 Data File : BF138922.D
 Acq On : 10 Aug 2024 21:25
 Operator : RC/JU
 Sample : P3467-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 919-J-WS-080224

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_F\Methods\8270-BF073024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138922.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.128	3	6	13	rVB	1780875	2686022	100.00%	23.872%
2	3.387	214	220	250	rBV	16722	52810	1.97%	0.469%
3	5.063	500	505	520	rVB	43381	64117	2.39%	0.570%
4	5.475	570	575	598	rVB	539669	728560	27.12%	6.475%
5	6.487	742	747	766	rBV	338062	489280	18.22%	4.349%
6	6.675	772	779	791	rBV	24248	39482	1.47%	0.351%
7	6.840	802	807	814	rBV	200084	244548	9.10%	2.173%
8	7.404	897	903	914	rBV	699769	941747	35.06%	8.370%
9	8.122	1019	1025	1041	rVB	242722	318194	11.85%	2.828%
10	8.434	1073	1078	1091	rBV	86250	118828	4.42%	1.056%
11	9.192	1201	1207	1212	rBV	1285790	1597697	59.48%	14.200%
12	9.869	1317	1322	1332	rBV	272251	343239	12.78%	3.051%
13	9.963	1332	1338	1344	rBV	103682	154239	5.74%	1.371%
14	10.663	1452	1457	1471	rBV	680225	886852	33.02%	7.882%
15	11.304	1562	1566	1570	rBV	31429	35621	1.33%	0.317%
16	11.357	1570	1575	1588	rVB	252655	323290	12.04%	2.873%
17	11.574	1606	1612	1620	rBV2	15479	28257	1.05%	0.251%
18	11.880	1659	1664	1679	rBV2	37610	68793	2.56%	0.611%
19	12.392	1746	1751	1757	rBV	28038	46159	1.72%	0.410%
20	12.751	1808	1812	1818	rVB	25906	30220	1.13%	0.269%
21	12.939	1838	1844	1849	rBV	878390	1109989	41.32%	9.865%
22	13.827	1990	1995	2005	rBV	38425	62328	2.32%	0.554%
23	13.998	2019	2024	2032	rVB	150318	194796	7.25%	1.731%
24	14.821	2159	2164	2167	rBV	41203	64126	2.39%	0.570%
25	14.868	2168	2172	2181	rVV5	37078	108981	4.06%	0.969%
26	14.980	2182	2191	2205	rVB3	49634	187725	6.99%	1.668%
27	15.462	2267	2273	2284	rVB	139578	216846	8.07%	1.927%
28	16.309	2410	2417	2427	rBV3	45438	108788	4.05%	0.967%

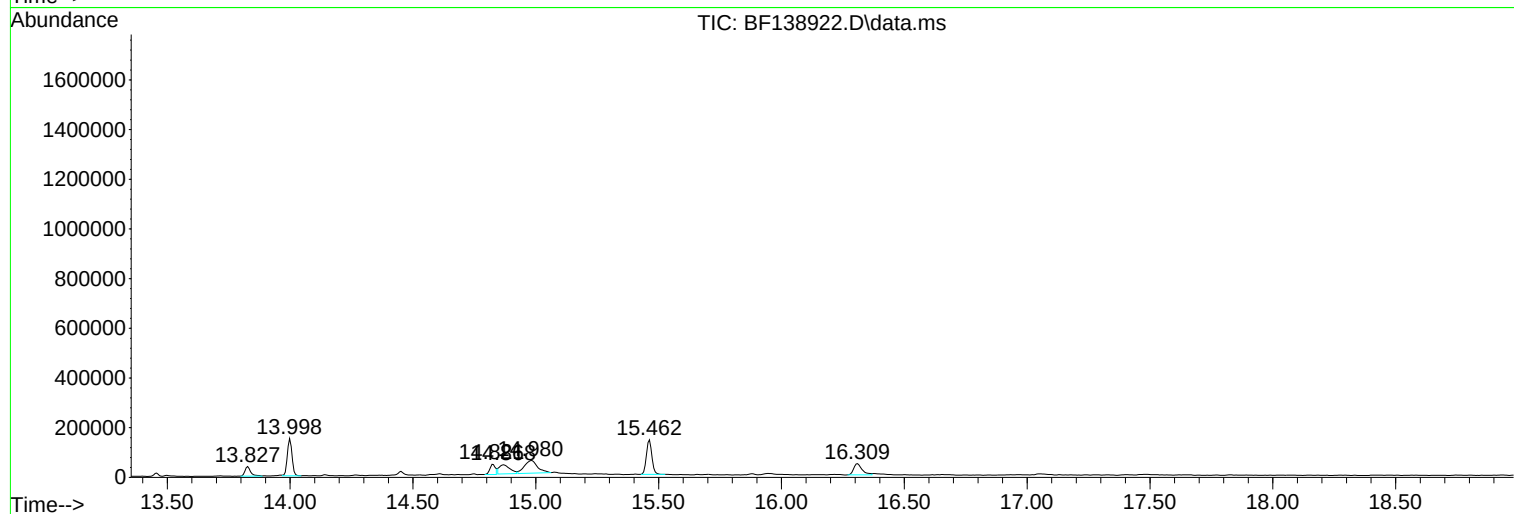
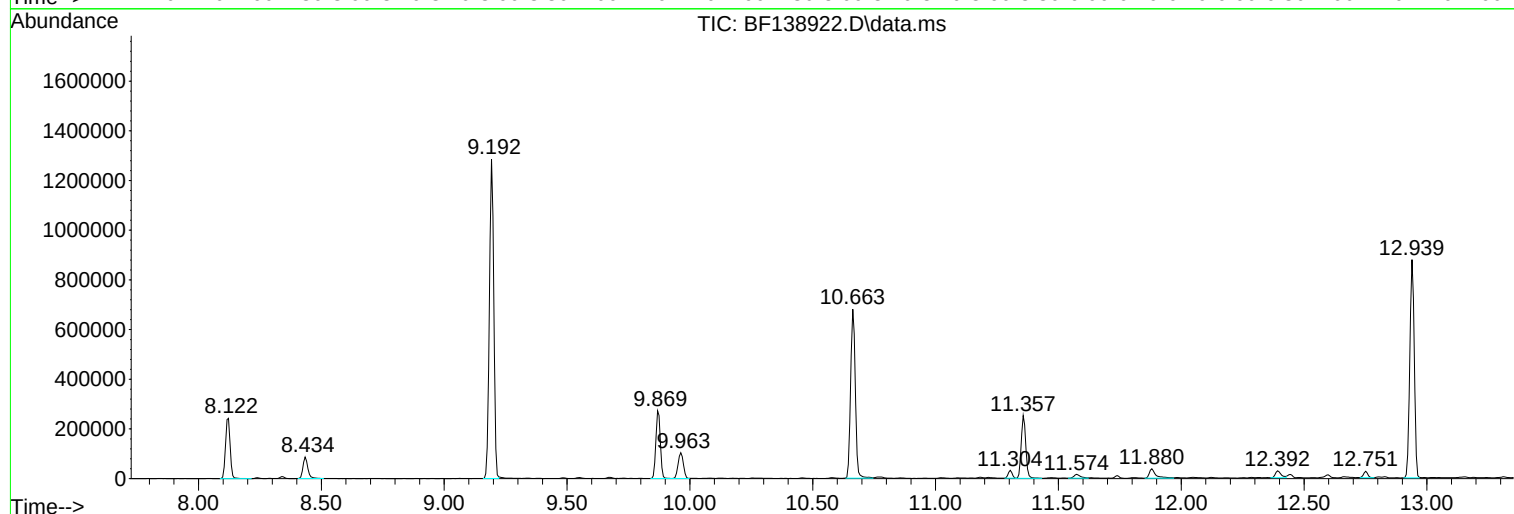
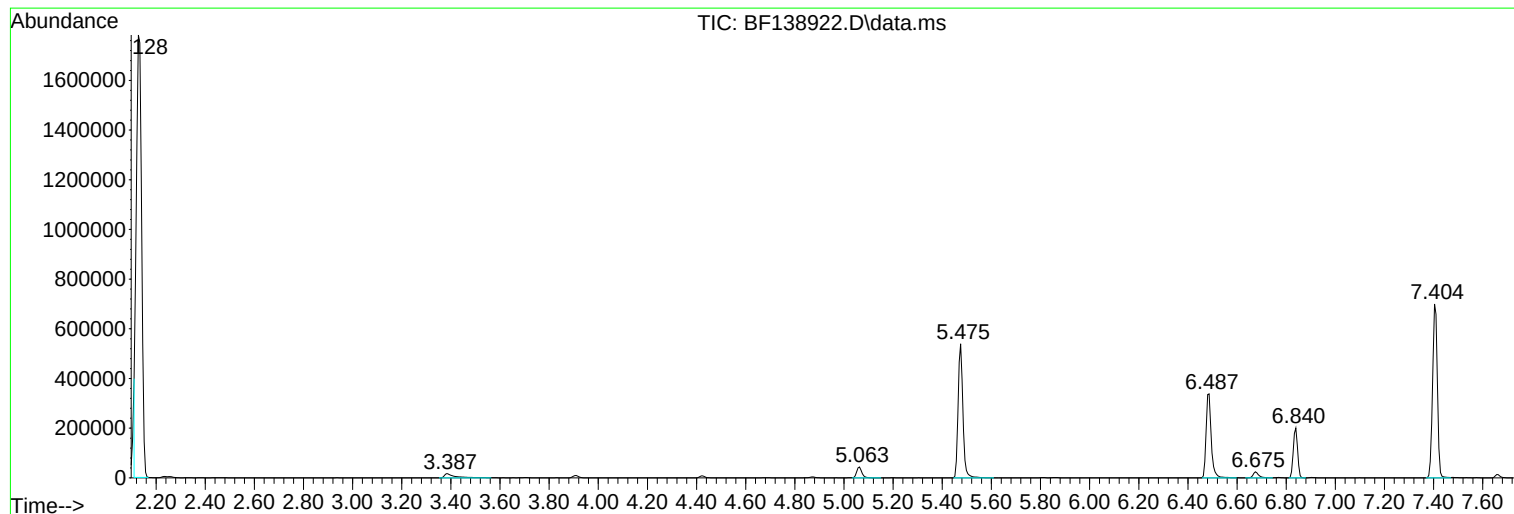
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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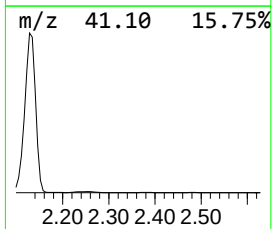
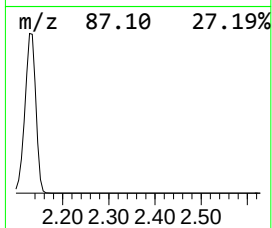
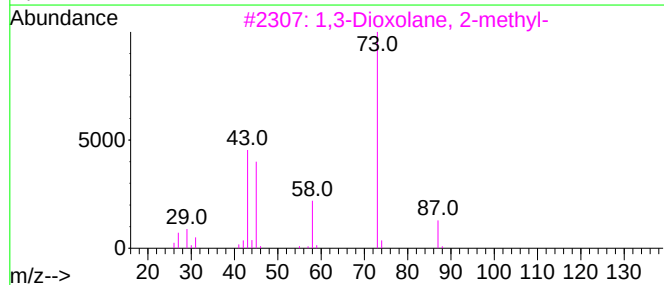
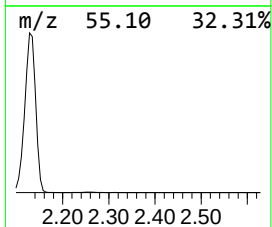
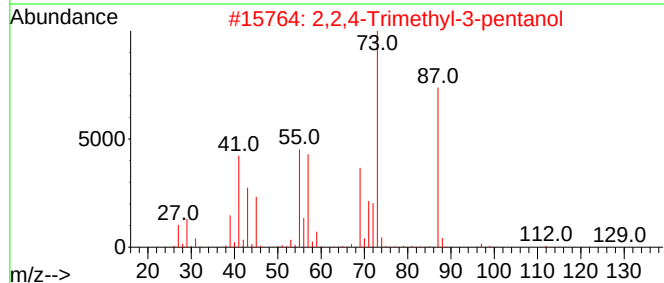
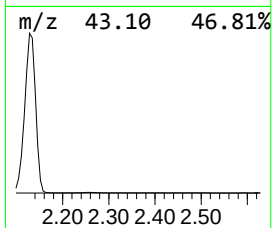
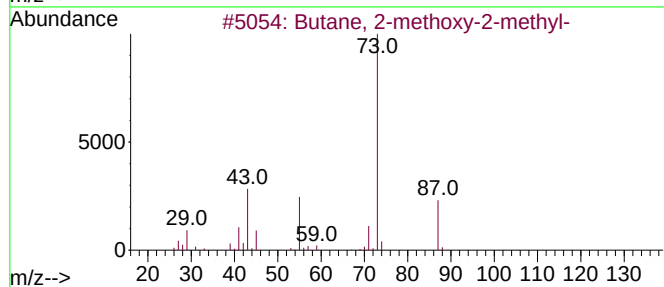
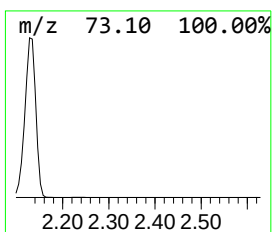
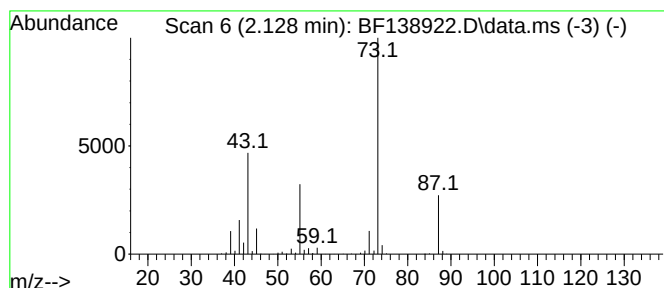
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.128	219.67 ng	2686020	1,4-Dichlorobenzene-d4	6.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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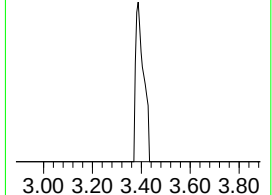
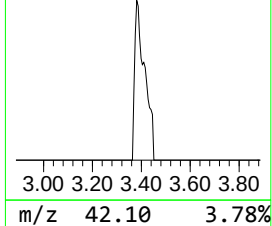
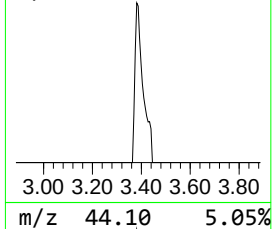
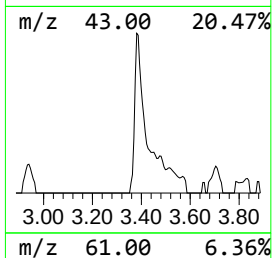
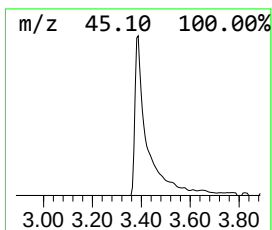
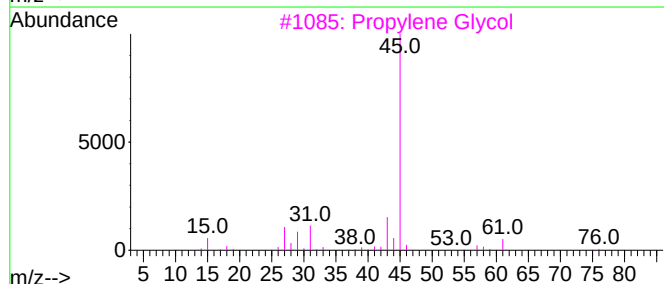
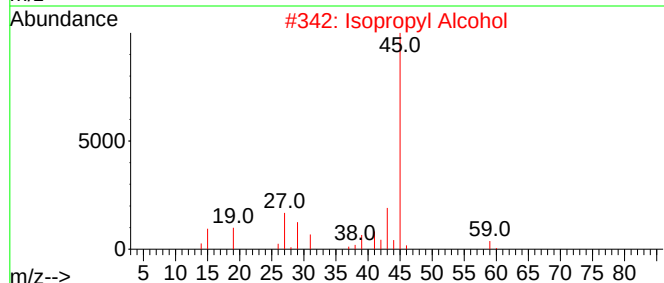
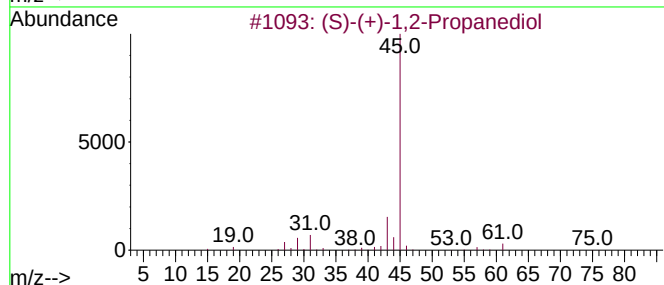
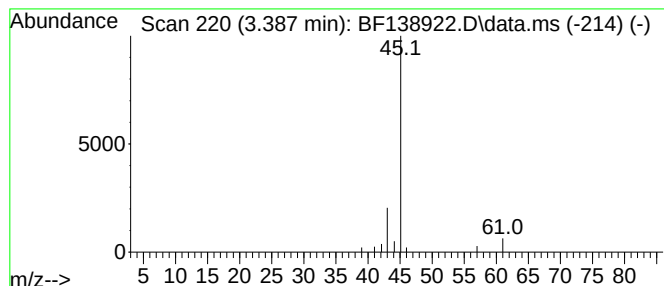
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 (S)-(+)-1,2-Propanediol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.387	4.32 ng	52810	1,4-Dichlorobenzene-d4	6.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	64
2	Isopropyl Alcohol			60	C3H8O	000067-63-0	9
3	Propylene Glycol			76	C3H8O2	000057-55-6	9
4	Formamide			45	CH3NO	000075-12-7	5
5	Propanoic acid, 2-hydroxy-, meth...			104	C4H8O3	000547-64-8	4



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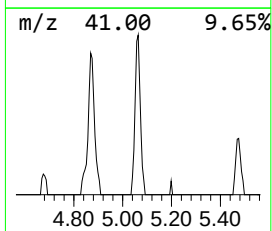
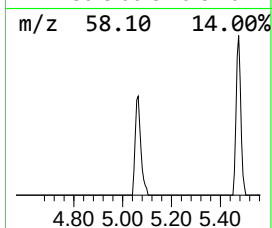
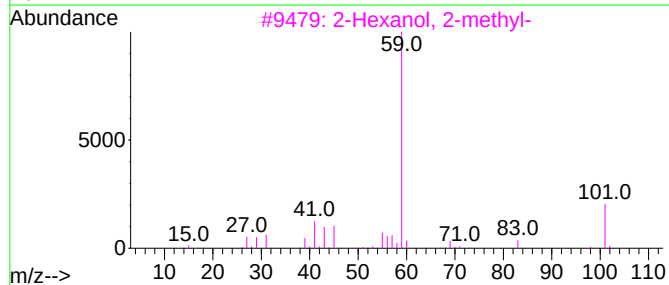
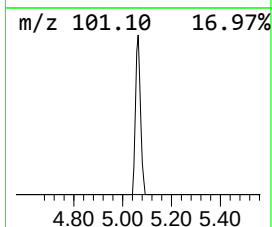
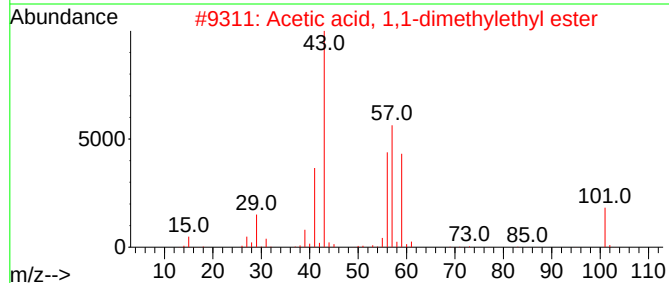
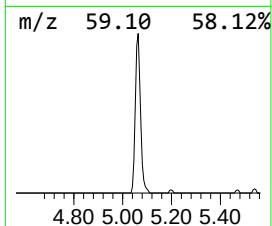
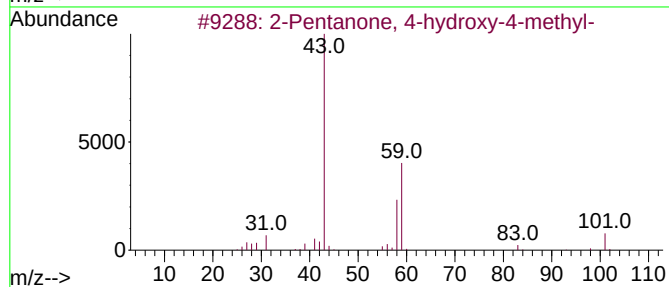
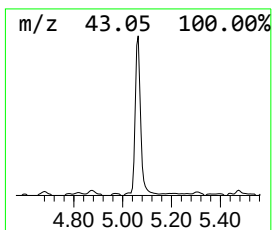
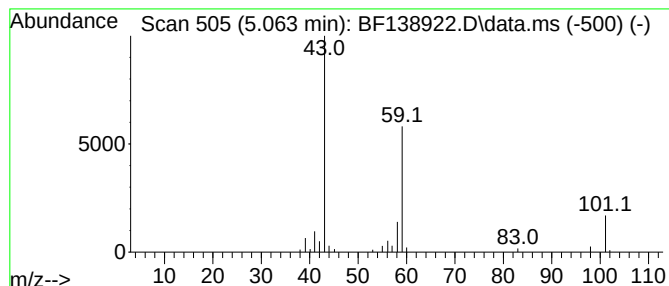
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.063	5.24 ng	64117	1,4-Dichlorobenzene-d4	6.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
4			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
5			3-Octanol	130	C8H18O	000589-98-0	33



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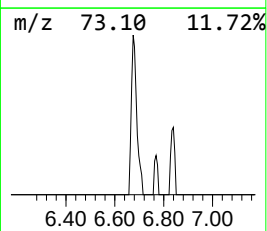
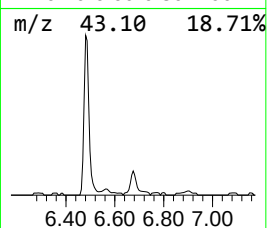
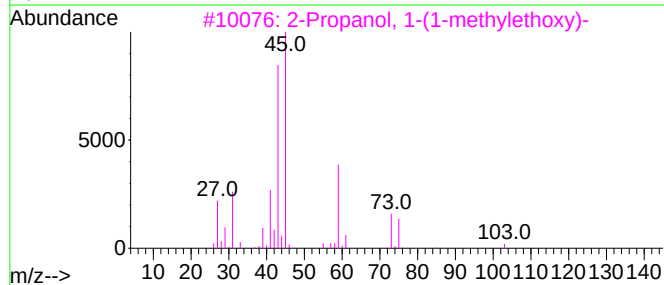
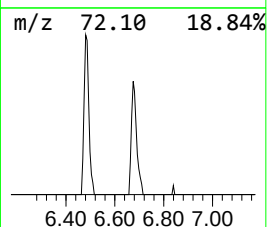
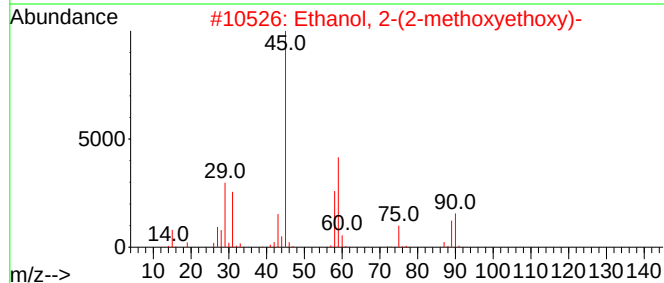
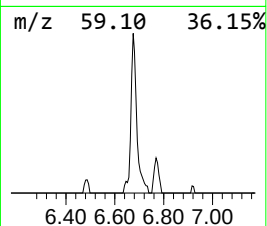
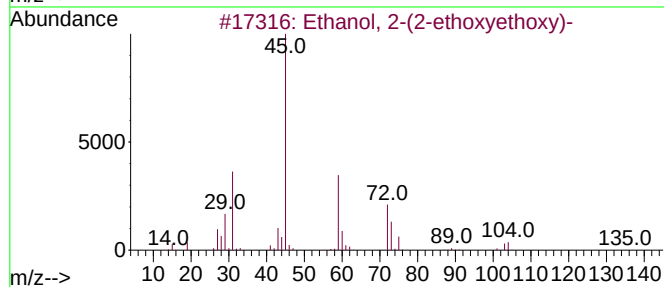
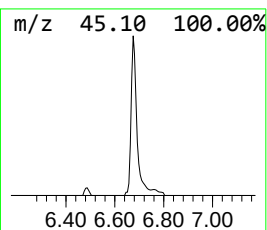
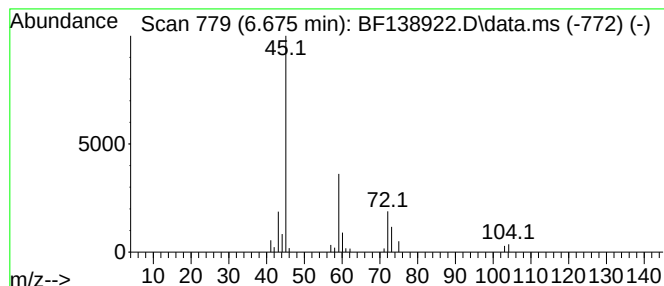
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.675	3.23 ng	39482	1,4-Dichlorobenzene-d4	6.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
2			Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	50
3			2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	45
4			2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	38
5			2-Methoxyisopropyl cyanoacetate	157	C7H11NO3	032804-79-8	33



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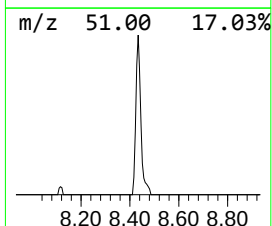
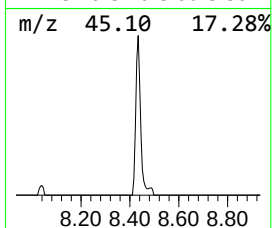
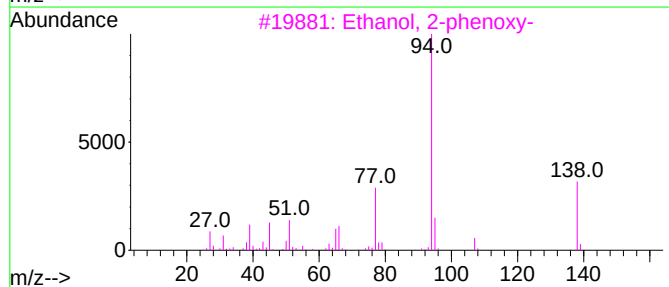
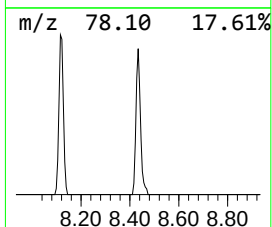
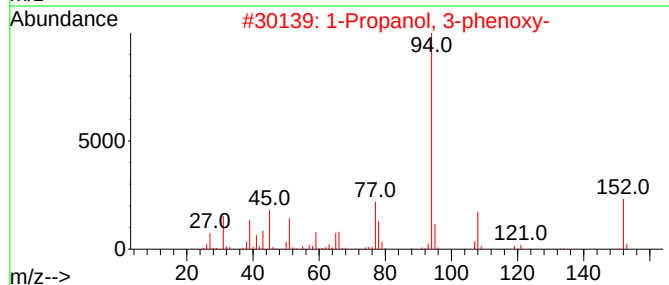
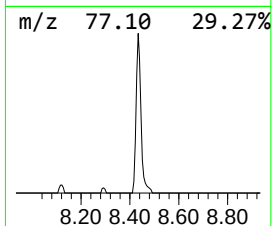
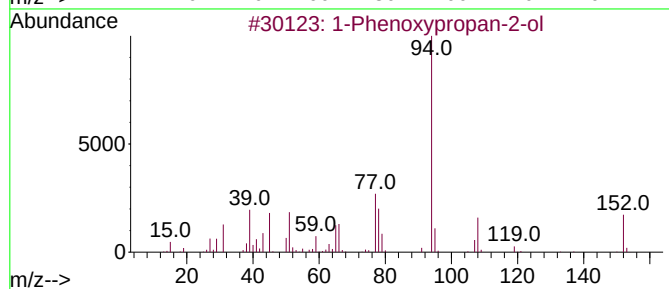
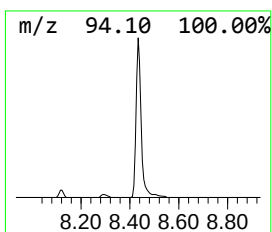
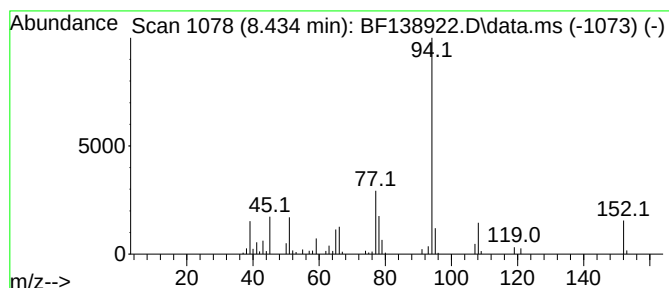
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.434	7.47 ng	118828	Naphthalene-d8	8.122

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
3			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	53
5			Benzene, (2-methylpropoxy)-	150	C10H14O	001126-75-6	53



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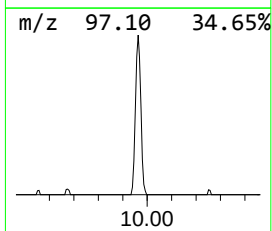
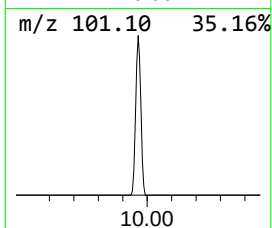
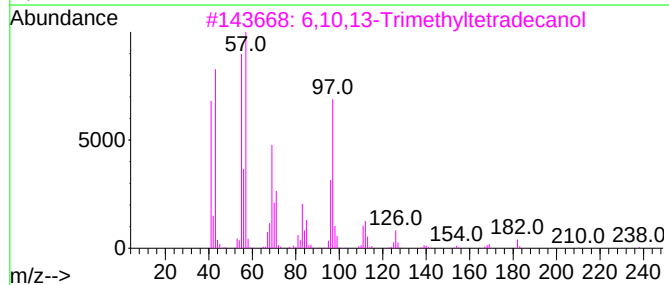
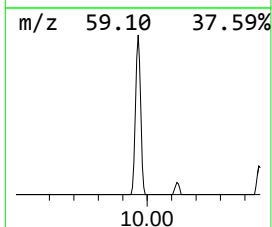
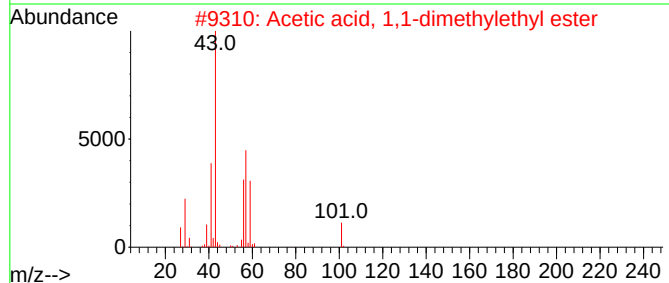
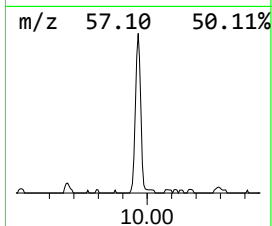
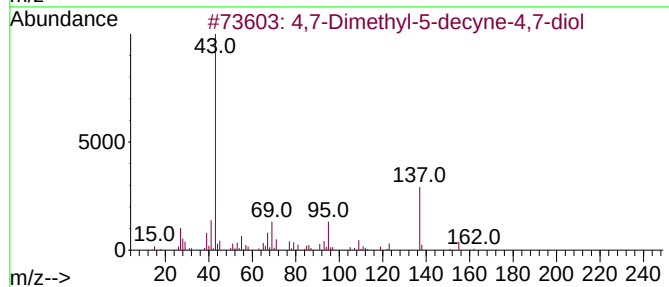
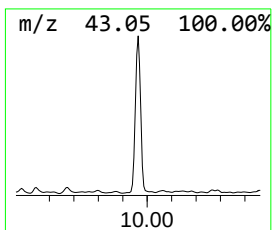
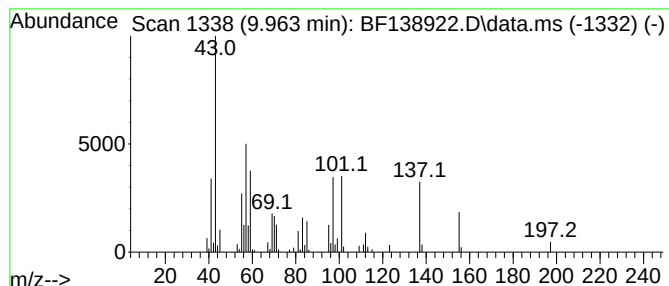
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown9.963 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.963	8.99 ng	154239	Acenaphthene-d10	9.869

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4,7-Dimethyl-5-decyne-4,7-diol	198	C12H22O2	000126-87-4	12
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	10
3		6,10,13-Trimethyltetradecanol	256	C17H36O	1000131-71-0	10
4		Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	10
5		2-Propionyloxytetradecane	270	C17H34O2	1000245-62-7	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
Sample : P3467-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
919-J-WS-080224

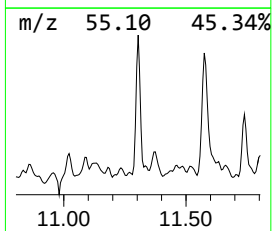
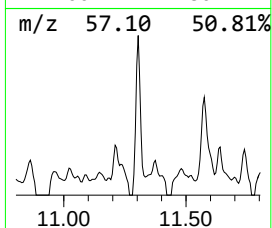
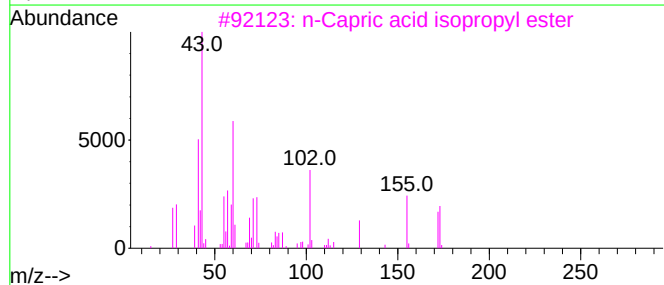
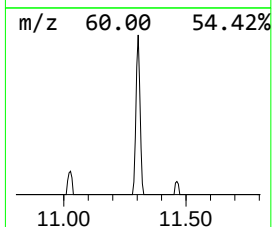
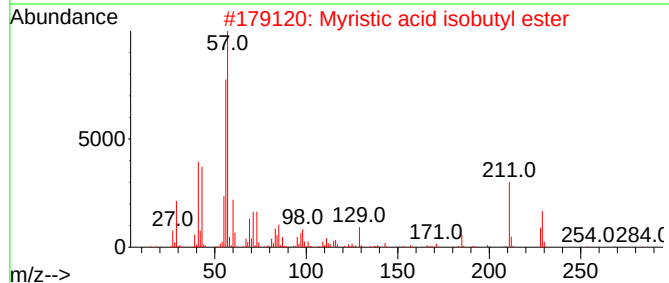
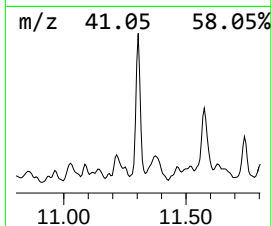
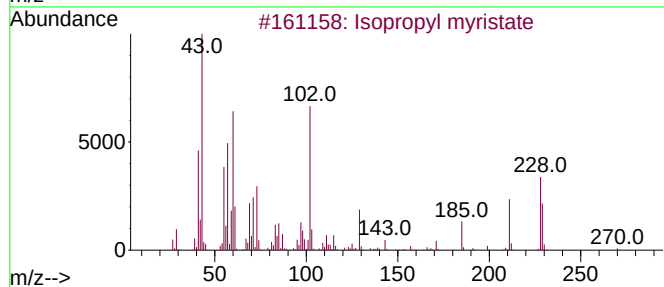
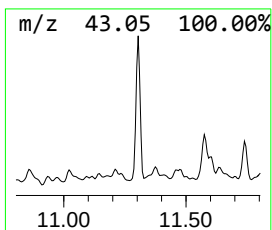
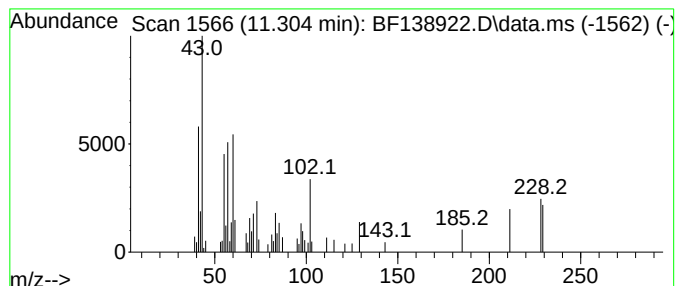
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Isopropyl myristate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.304	2.20 ng	35621	Phenanthrene-d10	11.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl myristate	270	C17H34O2	000110-27-0	90
2			Myristic acid isobutyl ester	284	C18H36O2	025263-97-2	47
3			n-Capric acid isopropyl ester	214	C13H26O2	002311-59-3	43
4			8-Methylnonanoic acid, isopropyl...	214	C13H26O2	1000452-00-6	38
5			Nonanoic acid, 1-methylethyl ester	200	C12H24O2	028267-32-5	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
Sample : P3467-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

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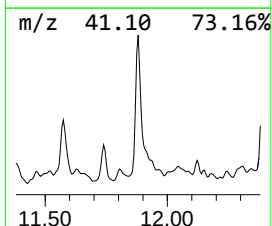
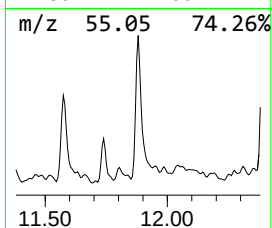
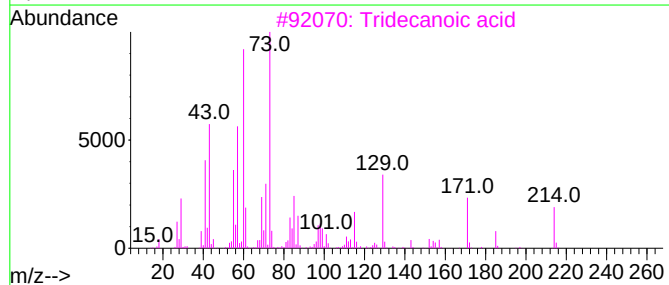
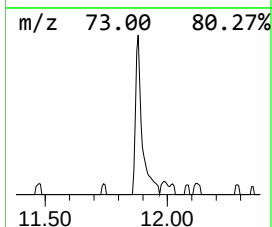
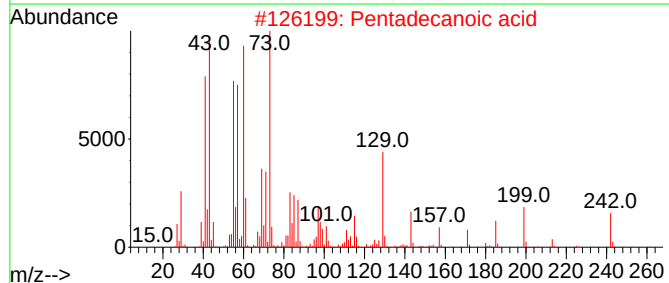
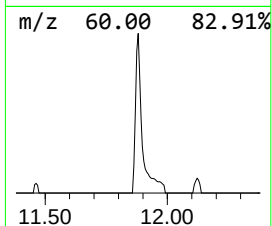
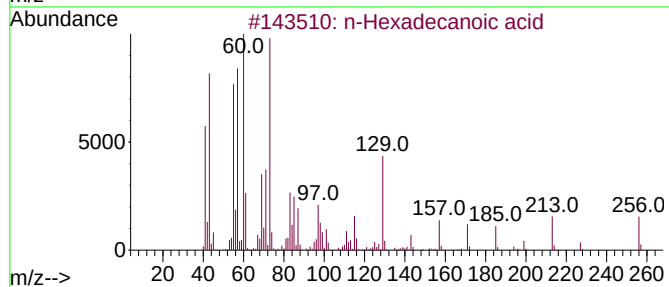
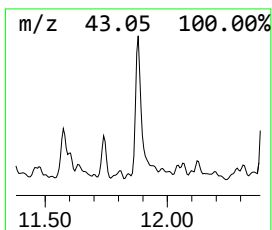
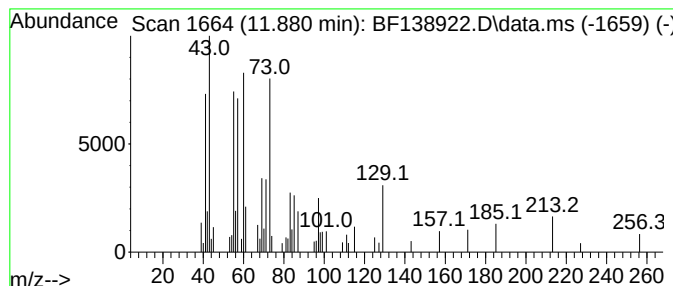
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.880	4.26 ng	68793	Phenanthrene-d10	11.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	74
3			Tridecanoic acid	214	C13H26O2	000638-53-9	72
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	53
5			Undecanoic acid	186	C11H22O2	000112-37-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
Sample : P3467-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

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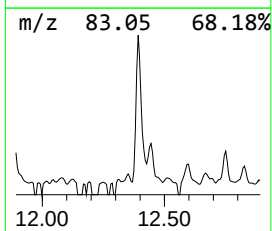
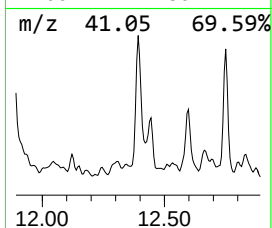
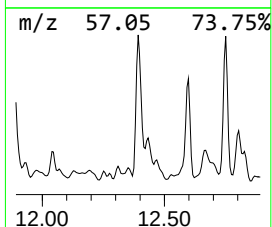
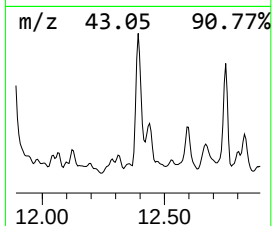
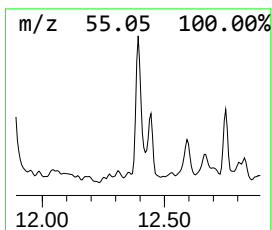
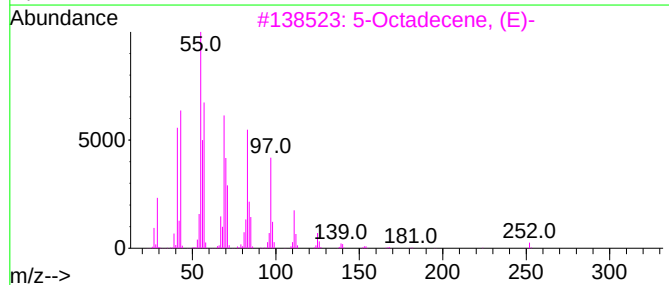
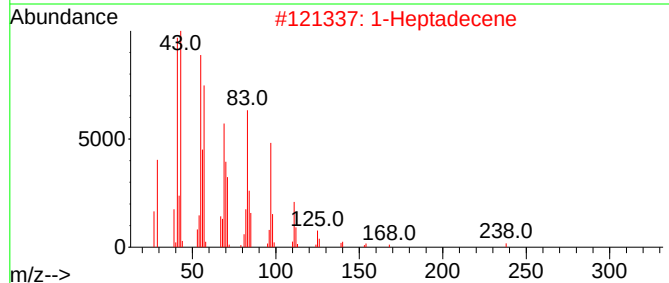
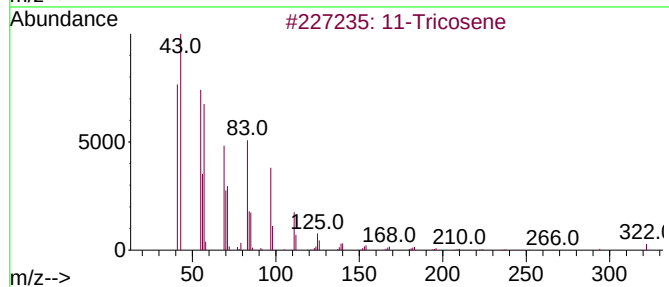
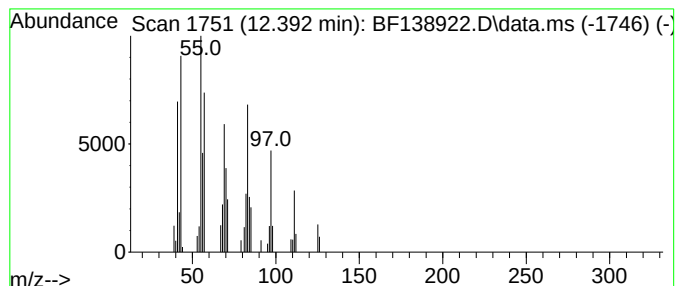
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 11-Tricosene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.392	2.86 ng	46159	Phenanthrene-d10	11.357

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		11-Tricosene	322	C23H46	052078-56-5	90
2		1-Heptadecene	238	C17H34	006765-39-5	87
3		5-Octadecene, (E)-	252	C18H36	007206-21-5	87
4		Cetene	224	C16H32	000629-73-2	86
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
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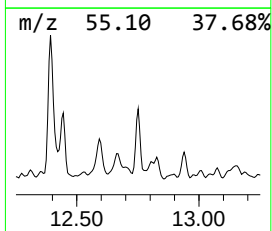
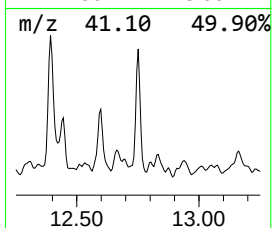
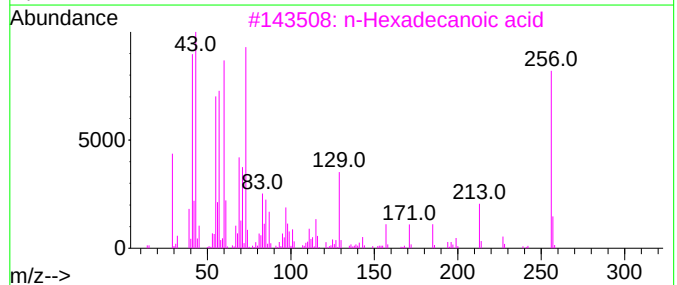
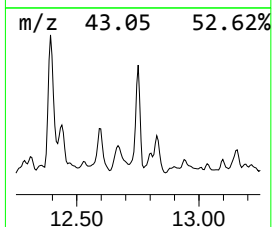
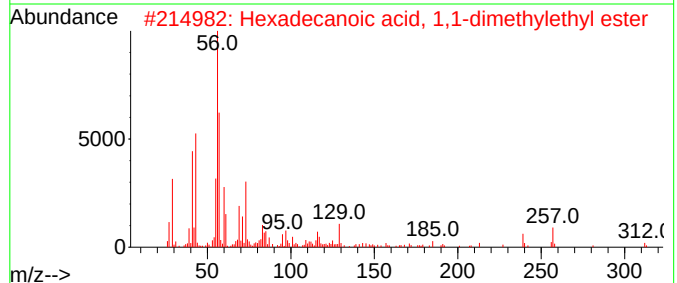
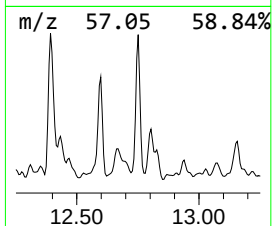
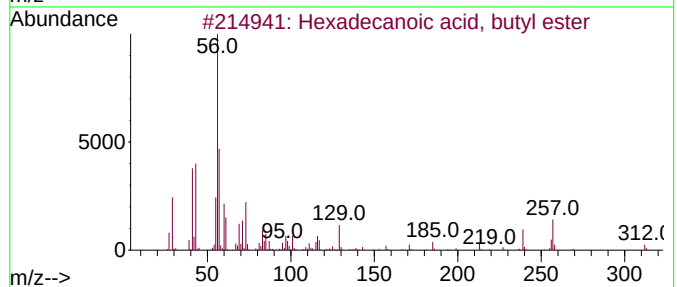
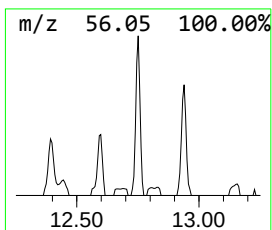
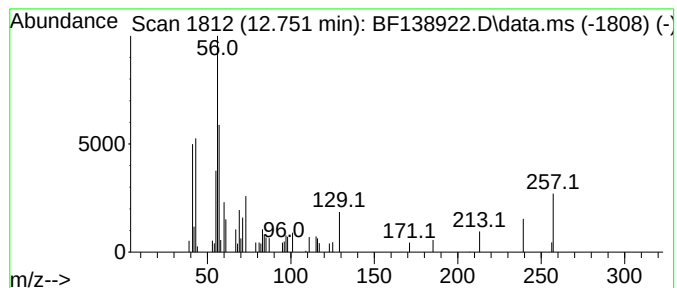
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecanoic acid, butyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.751	3.10 ng	30220	Chrysene-d12	13.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	87
2			Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	64
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	27
4			1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	27
5			1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
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Misc :
ALS Vial : 23 Sample Multiplier: 1

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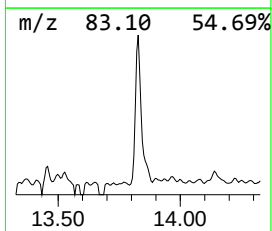
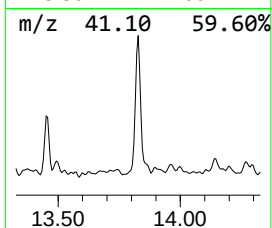
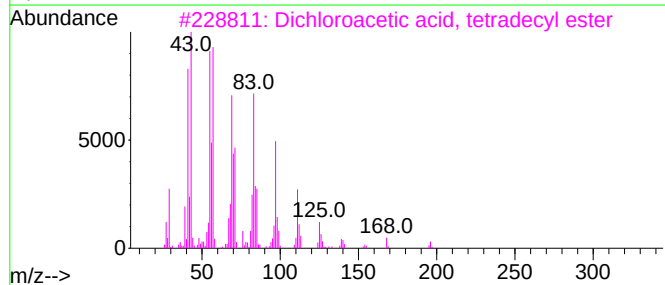
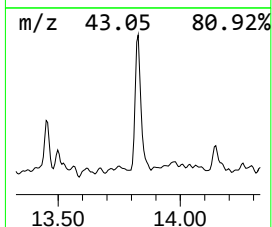
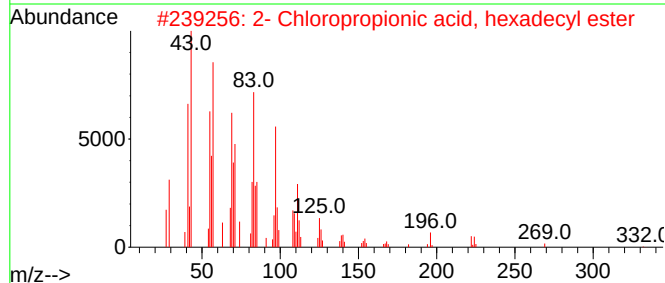
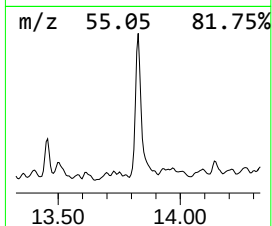
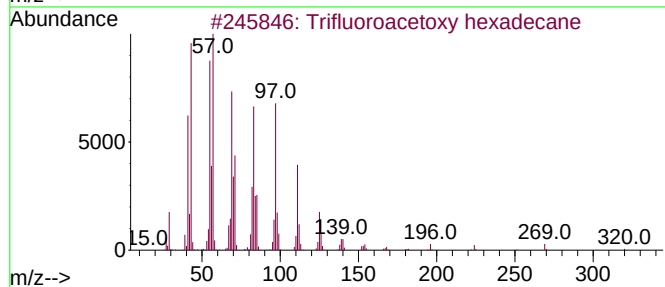
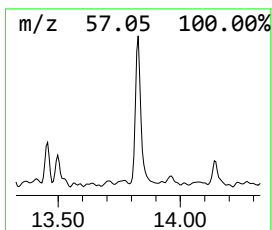
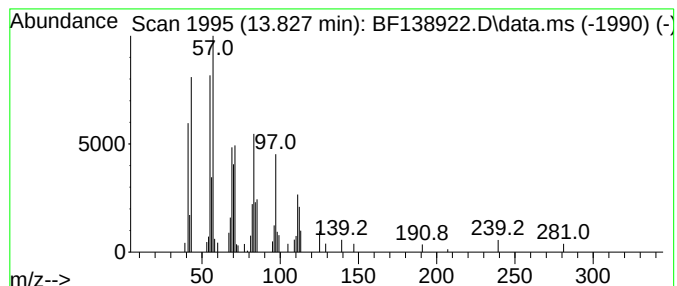
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Trifluoroacetoxy hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.827	6.40 ng	62328	Chrysene-d12	13.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	87
2			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	87
3			Dichloroacetic acid, tetradecyl ...	324	C16H30Cl2O2	083005-02-1	87
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	87
5			1-Docosene	308	C22H44	001599-67-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
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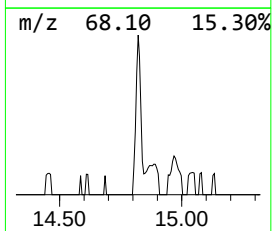
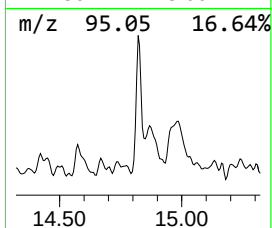
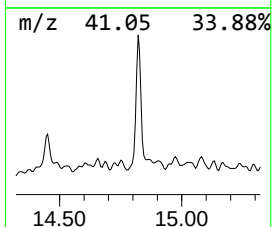
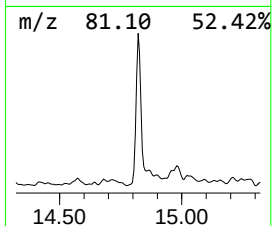
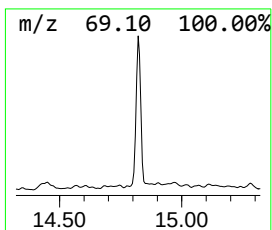
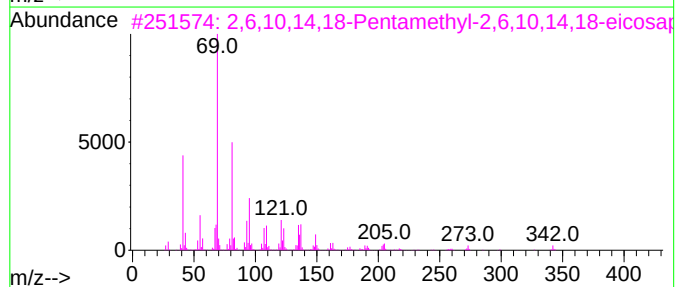
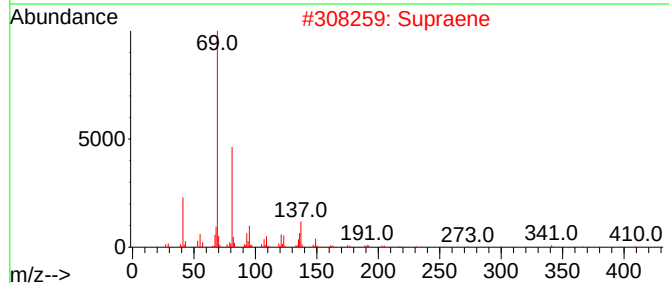
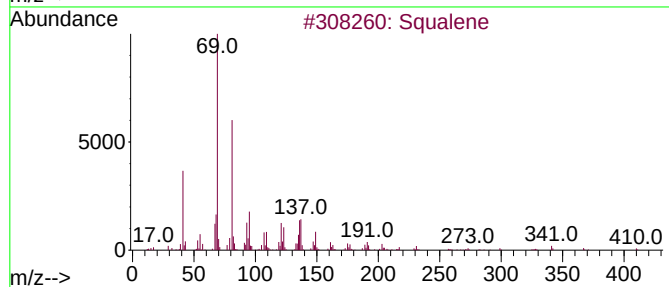
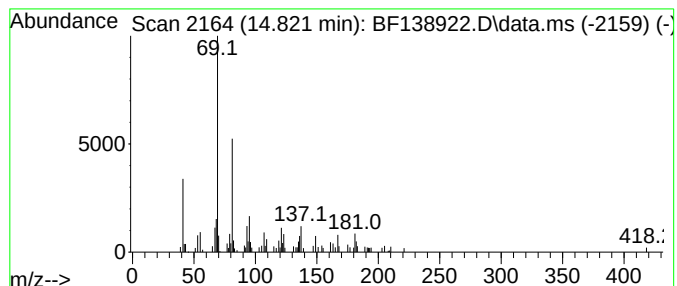
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Squalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.821	5.91 ng	64126	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	86
2			Supraene	410	C30H50	007683-64-9	80
3			2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42	075581-03-2	80
4			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
5			Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
Sample : P3467-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
919-J-WS-080224

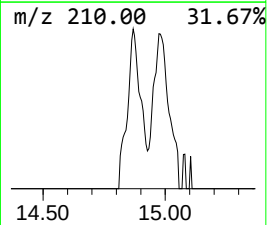
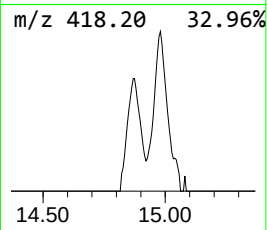
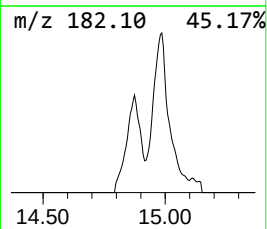
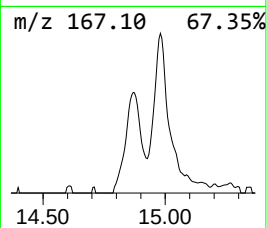
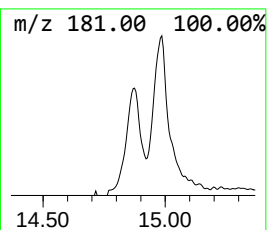
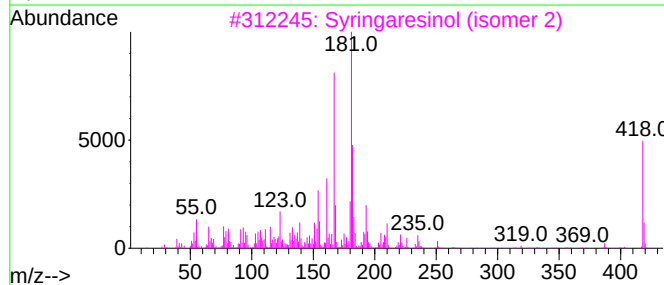
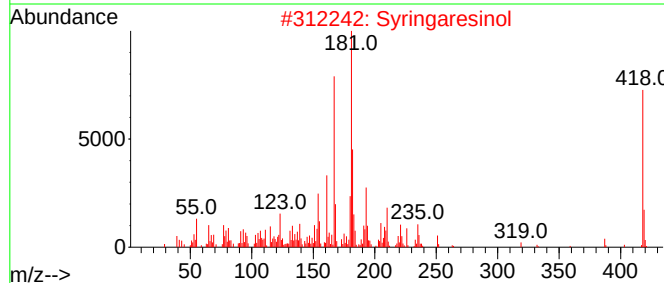
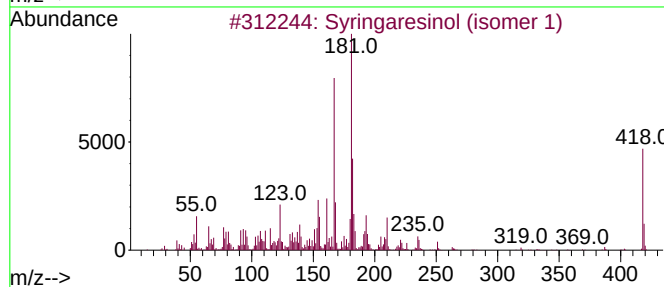
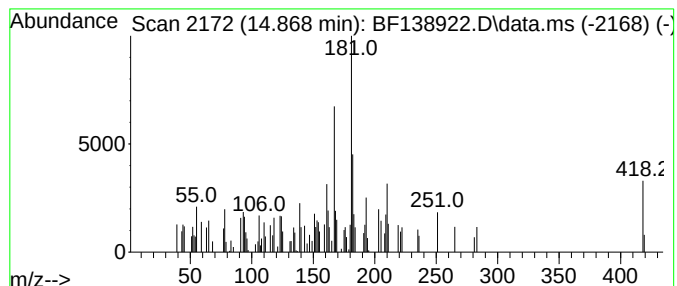
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Syringaresinol (isomer 1) Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.868	10.05 ng	108981	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Syringaresinol (isomer 1)	418	C22H26O8	1000495-97-0	98
2			Syringaresinol	418	C22H26O8	000487-35-4	93
3			Syringaresinol (isomer 2)	418	C22H26O8	1000495-97-1	90
4			1,2,3,3a,4,5,6,10b-Octahydroflu...	210	C16H18	1010080-20-5	58
5			1,10-phenanthroline-2(1H)-one, 1-...	210	C13H10N2O	1000400-61-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
Sample : P3467-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
919-J-WS-080224

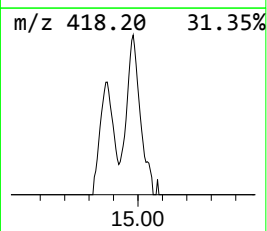
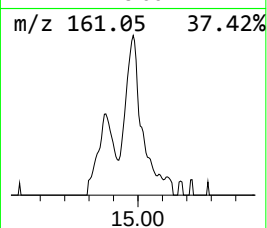
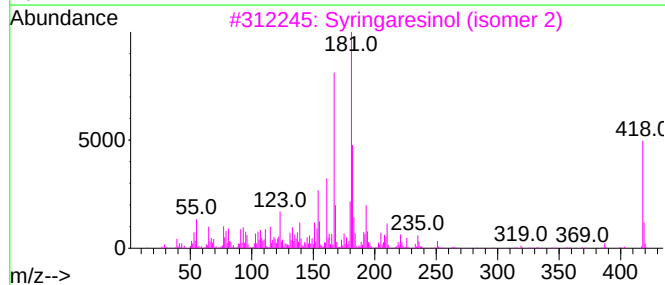
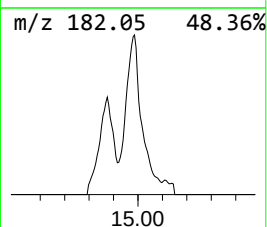
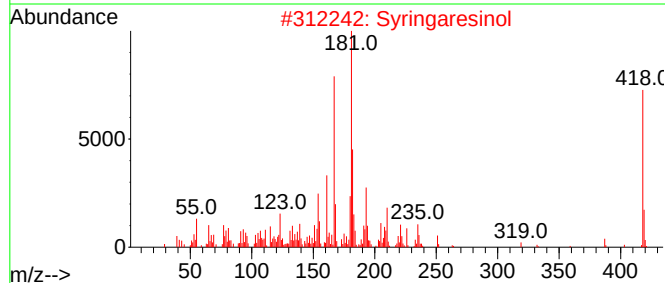
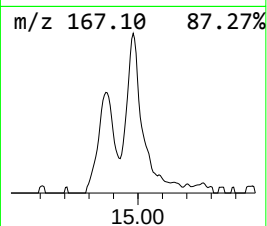
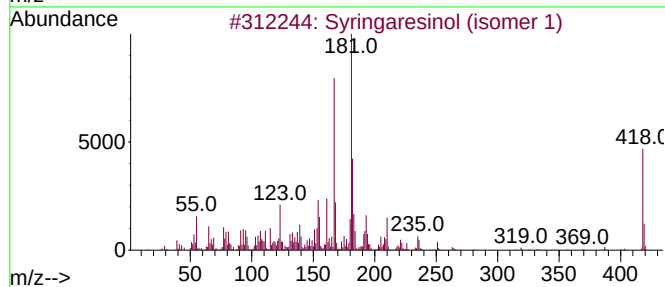
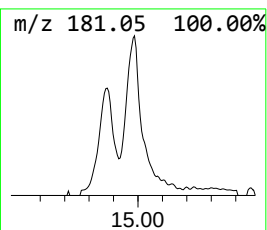
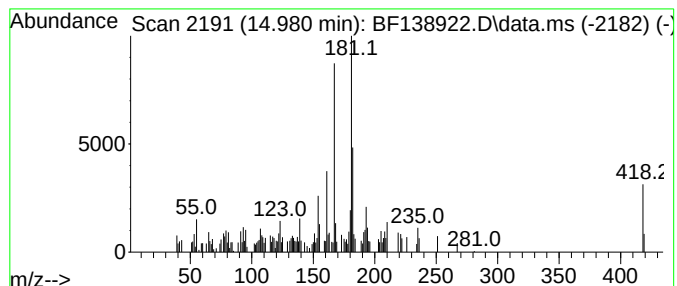
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Syringaresinol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.980	17.31 ng	187725	Perylene-d12	15.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Syringaresinol (isomer 1)	418	C22H26O8	1000495-97-0	94
2		Syringaresinol	418	C22H26O8	000487-35-4	93
3		Syringaresinol (isomer 2)	418	C22H26O8	1000495-97-1	55
4		2,6-Pyridinedimethanol, 4-(dimet...	182	C9H14N2O2	001882-25-3	38
5		phosphorane, (3,5-dimethylphenyl...	182	C10H15OP	1000404-87-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
Sample : P3467-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
919-J-WS-080224

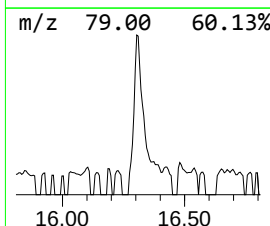
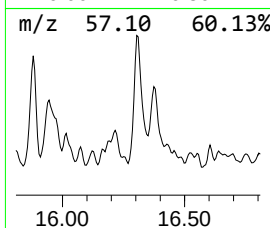
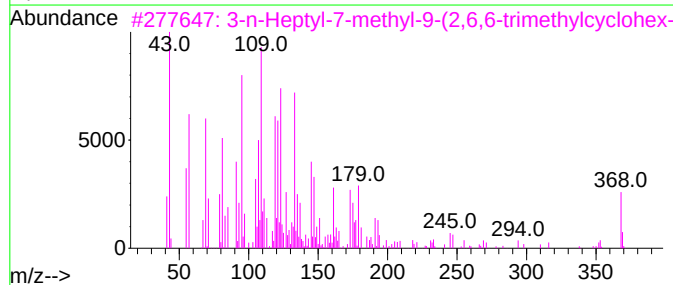
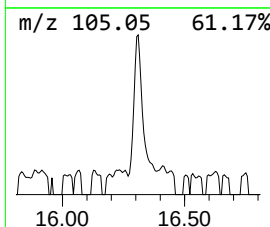
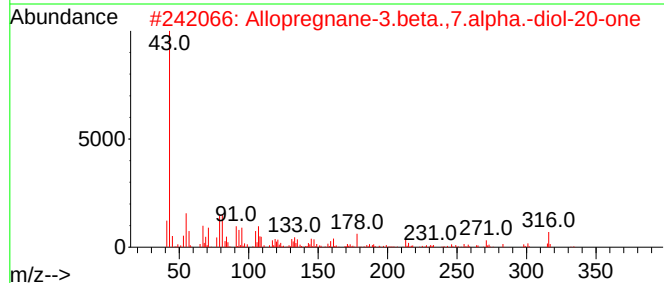
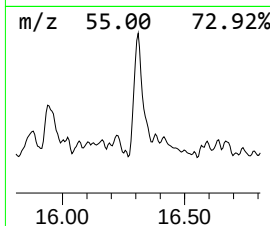
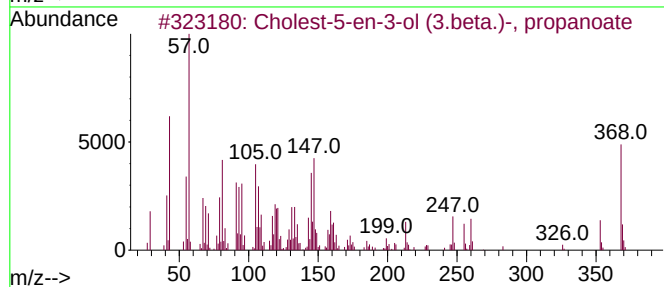
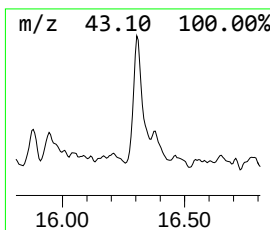
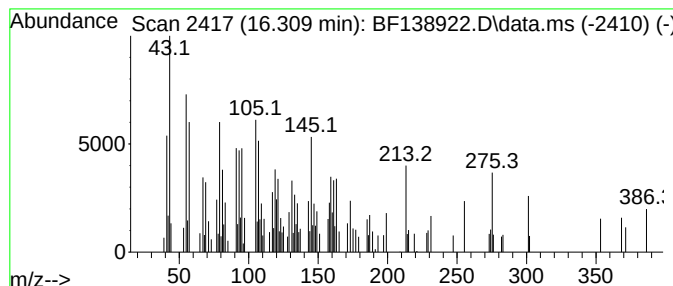
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown16.310 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.310	10.03 ng	108788	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholest-5-en-3-ol (3.beta.)-, pr...	442	C30H50O2	000633-31-8	38
2			Allopregnane-3.beta.,7.alpha.-di...	334	C21H34O3	000566-20-1	35
3			3-n-Heptyl-7-methyl-9-(2,6,6-tri...	368	C26H40O	1000216-09-3	35
4			Boscartol F	300	C20H28O2	1486443-17-7	25
5			Methyl 12-acetyl-7-desoxycholate	448	C27H44O5	055547-48-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081024\
Data File : BF138922.D
Acq On : 10 Aug 2024 21:25
Operator : RC/JU
Sample : P3467-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
919-J-WS-080224

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.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
Butane, 2-metho...	2.128	219.7	ng	2686020	1	6.840	244548	20.0
(S)-(+)-1,2-Pro...	3.387	4.3	ng	52810	1	6.840	244548	20.0
2-Pentanone, 4-...	5.063	5.2	ng	64117	1	6.840	244548	20.0
Ethanol, 2-(2-e...	6.675	3.2	ng	39482	1	6.840	244548	20.0
1-Phenoxypropan...	8.434	7.5	ng	118828	2	8.122	318194	20.0
unknown9.963	9.963	9.0	ng	154239	3	9.869	343239	20.0
Isopropyl myris...	11.304	2.2	ng	35621	4	11.357	323290	20.0
n-Hexadecanoic ...	11.880	4.3	ng	68793	4	11.357	323290	20.0
11-Tricosene	12.392	2.9	ng	46159	4	11.357	323290	20.0
Hexadecanoic ac...	12.751	3.1	ng	30220	5	13.998	194796	20.0
Trifluoroacetox...	13.827	6.4	ng	62328	5	13.998	194796	20.0
Squalene	14.821	5.9	ng	64126	6	15.463	216846	20.0
Syringaresinol ...	14.868	10.1	ng	108981	6	15.463	216846	20.0
Syringaresinol	14.980	17.3	ng	187725	6	15.463	216846	20.0
unknown16.310	16.310	10.0	ng	108788	6	15.463	216846	20.0