

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
 Data File : BF139007.D
 Acq On : 14 Aug 2024 21:47
 Operator : RC/JU
 Sample : P3475-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-866-N-S0-1.0-1.5-080524

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: BF139007.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	214	218	226	rBV	7701	20620	2.22%	0.330%
2	5.057	499	504	524	rVB	70799	111411	12.01%	1.784%
3	5.469	569	574	596	rBV	666923	927978	100.00%	14.857%
4	6.481	741	746	769	rBV	653350	906313	97.67%	14.511%
5	6.681	772	780	790	rBV	11872	23647	2.55%	0.379%
6	6.834	802	806	812	rVB	162397	198133	21.35%	3.172%
7	6.987	827	832	836	rBV	19218	23269	2.51%	0.373%
8	7.398	897	902	919	rBV	407522	541729	58.38%	8.673%
9	8.116	1019	1024	1039	rVB	170899	209677	22.60%	3.357%
10	8.434	1073	1078	1089	rBV	35903	55504	5.98%	0.889%
11	9.186	1201	1206	1216	rBV	558276	672672	72.49%	10.770%
12	9.869	1317	1322	1327	rBV	155167	194605	20.97%	3.116%
13	9.957	1332	1337	1344	rBV	27459	40803	4.40%	0.653%
14	10.657	1451	1456	1471	rBV	365116	470309	50.68%	7.530%
15	11.351	1569	1574	1586	rVB	170175	238722	25.72%	3.822%
16	11.875	1654	1663	1673	rBV7	12198	29409	3.17%	0.471%
17	12.569	1777	1781	1786	rBV	42550	51441	5.54%	0.824%
18	12.798	1814	1820	1827	rBV	34827	50163	5.41%	0.803%
19	12.933	1838	1843	1848	rBV	637150	791891	85.34%	12.679%
20	13.827	1988	1995	2008	rBV3	40676	93390	10.06%	1.495%
21	13.998	2019	2024	2033	rVB	140974	207679	22.38%	3.325%
22	14.463	2096	2103	2114	rBV3	22061	67087	7.23%	1.074%
23	14.733	2145	2149	2155	rBV2	21959	34100	3.67%	0.546%
24	14.892	2172	2176	2182	rVB4	13427	18078	1.95%	0.289%
25	15.039	2197	2201	2204	rBV	13290	22064	2.38%	0.353%
26	15.074	2204	2207	2211	rVB	15039	20332	2.19%	0.326%
27	15.121	2211	2215	2218	rBV4	6675	12246	1.32%	0.196%
28	15.339	2249	2252	2258	rVB	8696	13324	1.44%	0.213%
29	15.404	2259	2263	2267	rBV	10241	14927	1.61%	0.239%
30	15.463	2268	2273	2285	rVB	79633	125400	13.51%	2.008%
31	15.651	2301	2305	2310	rVB4	7172	11070	1.19%	0.177%
32	15.874	2338	2343	2349	rVB	12649	20949	2.26%	0.335%
33	16.668	2473	2478	2484	rVB6	6508	13603	1.47%	0.218%
34	16.951	2522	2526	2533	rVB5	7332	13329	1.44%	0.213%

Sum of corrected areas: 6245874

Instrument :

BNA_F

ClientSampleId :

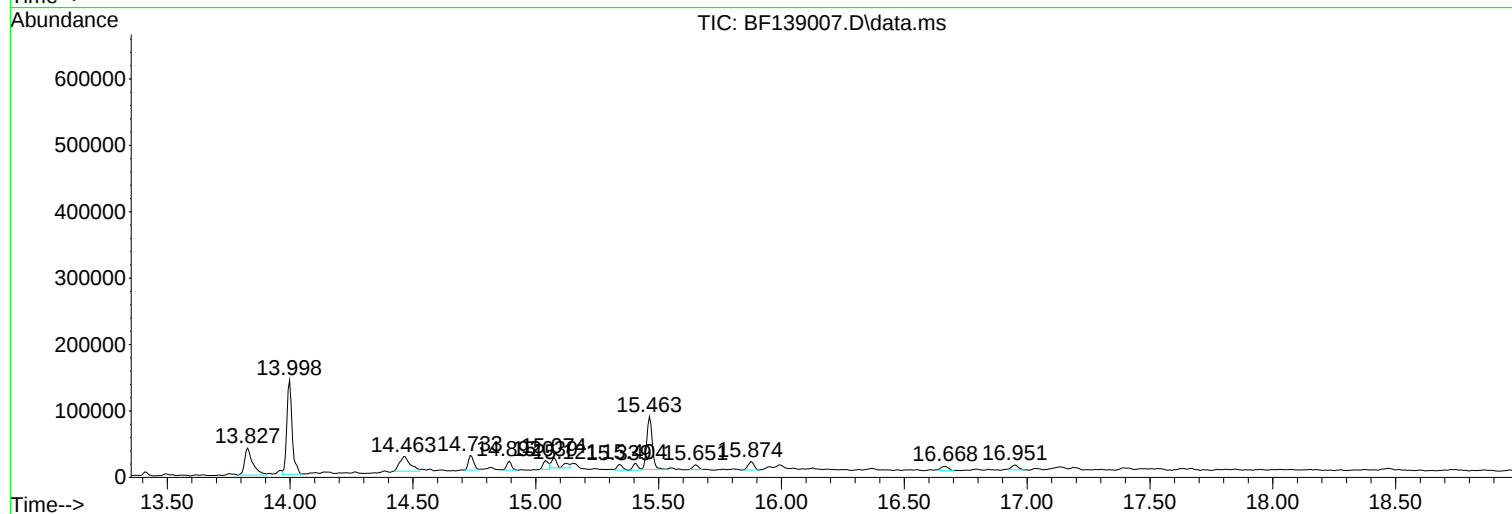
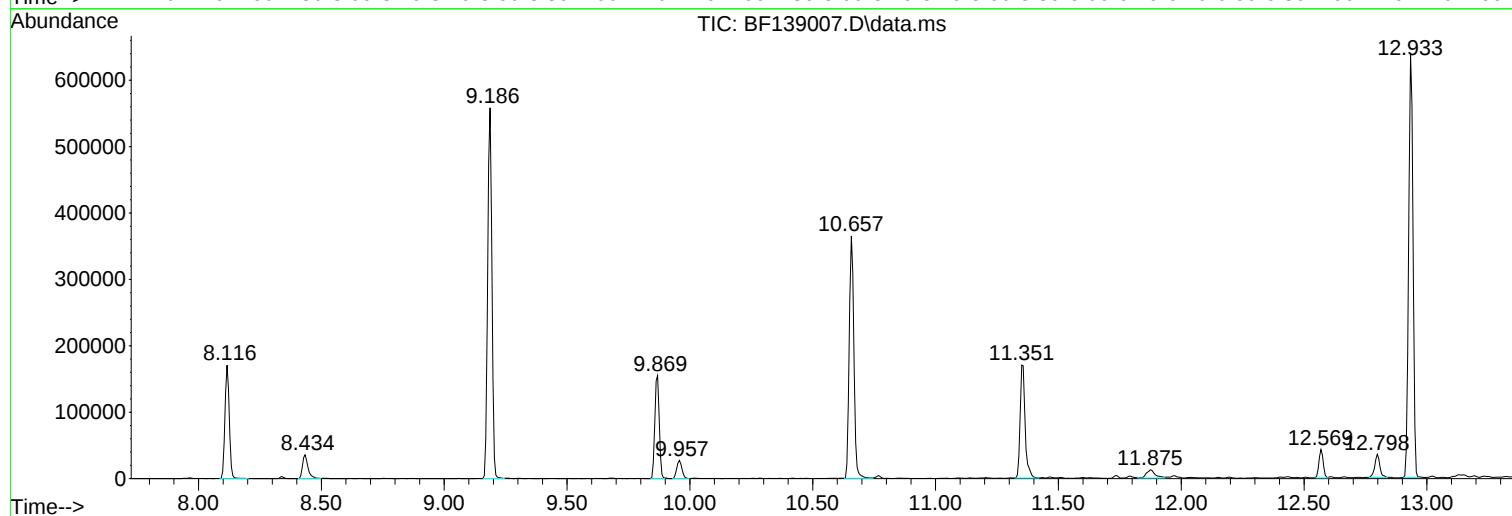
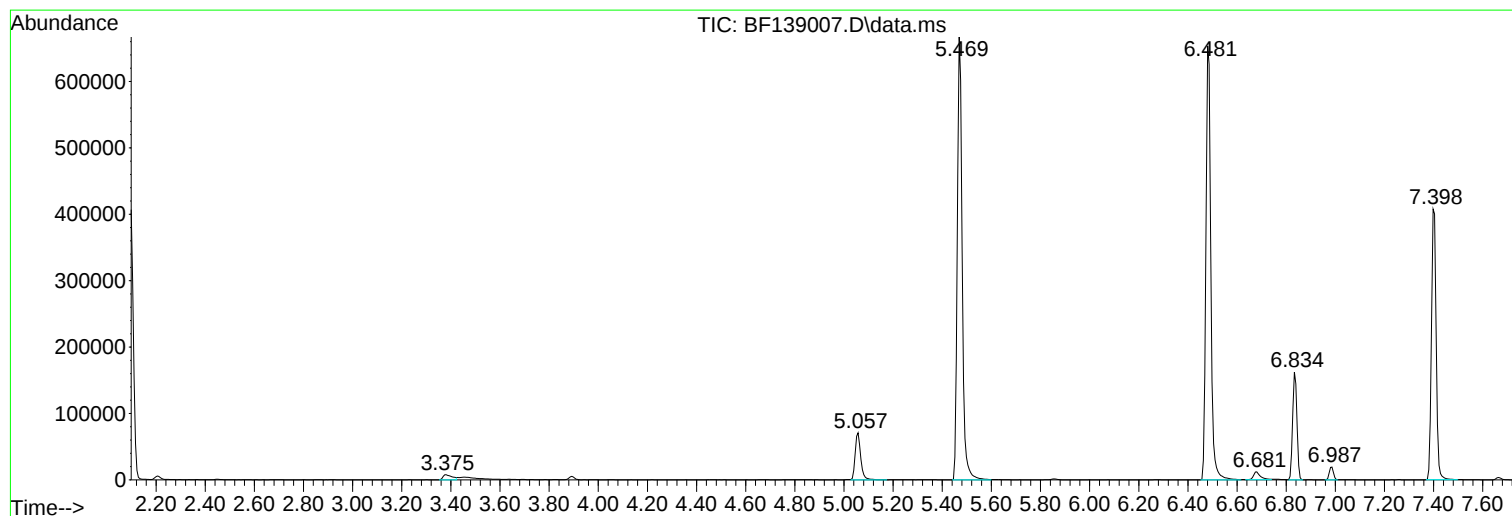
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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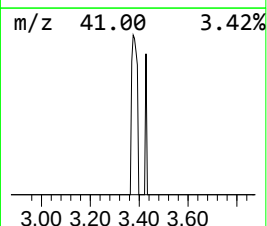
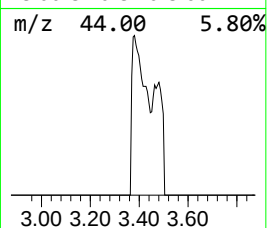
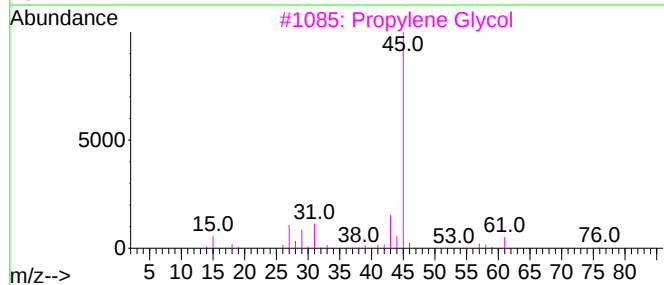
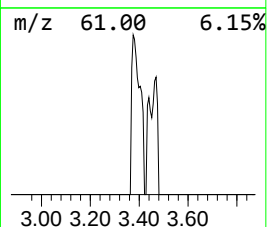
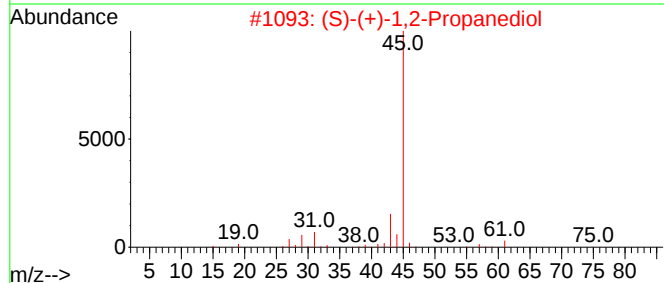
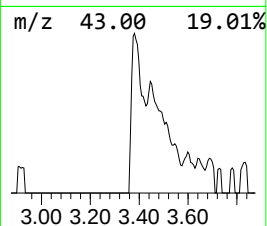
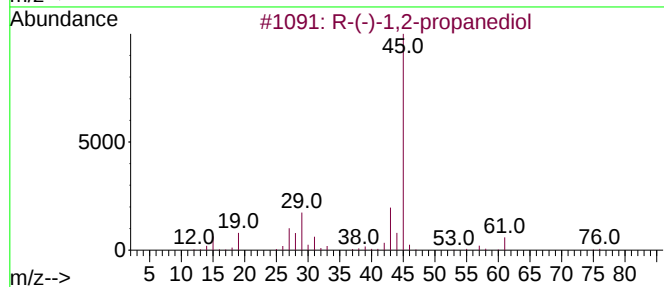
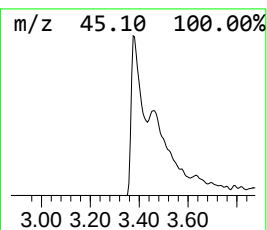
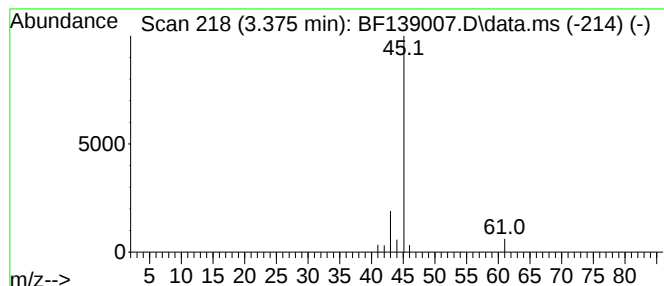
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 R-(-)-1,2-propanediol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.375	2.08 ng	20620	1,4-Dichlorobenzene-d4	6.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	64
2	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	9
3	Propylene Glycol			76	C3H8O2	000057-55-6	9
4	Formamide			45	CH3NO	000075-12-7	7
5	Propanoic acid, 2-hydroxy-, meth...			104	C4H8O3	000547-64-8	4



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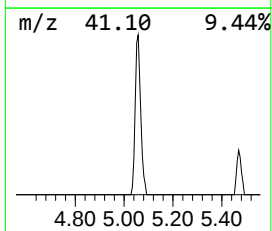
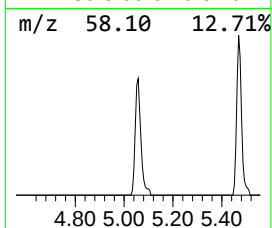
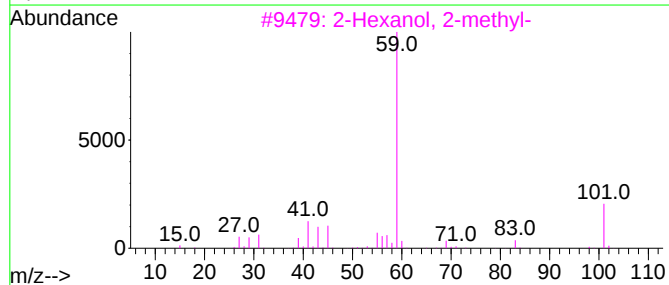
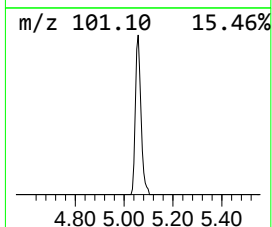
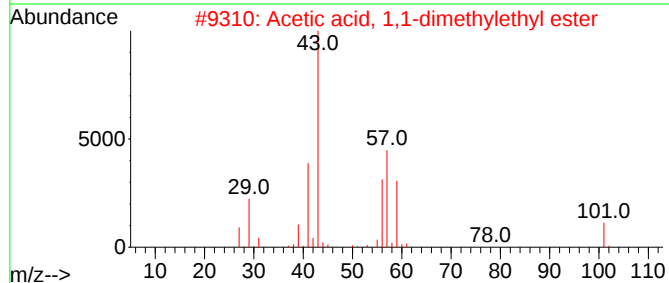
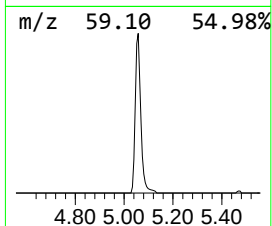
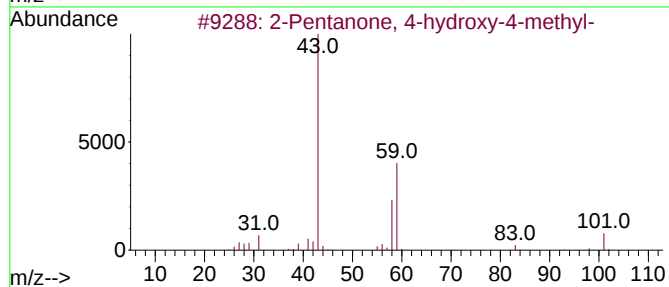
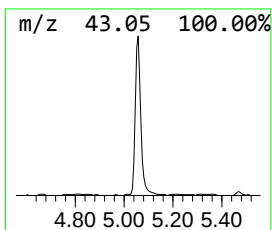
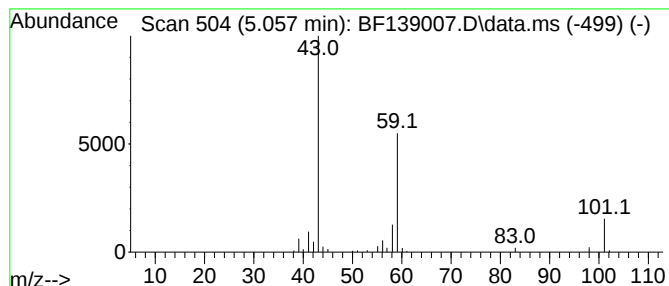
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.057	11.25 ng	111411	1,4-Dichlorobenzene-d4	6.834

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	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33		
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		



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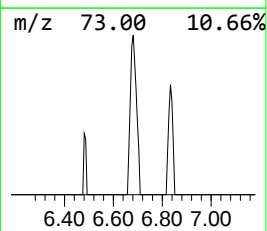
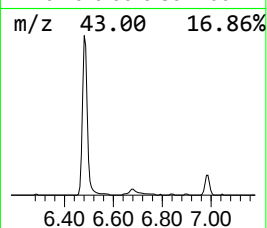
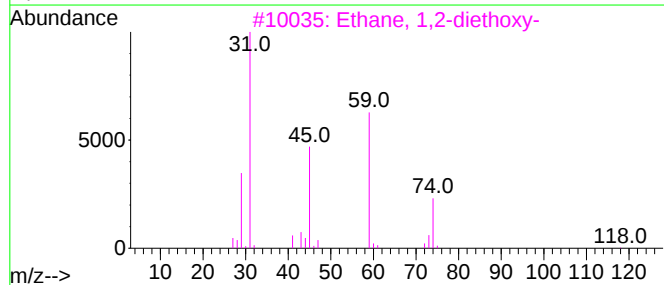
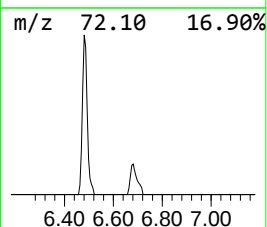
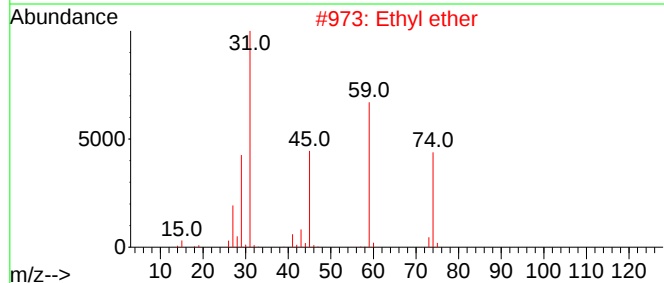
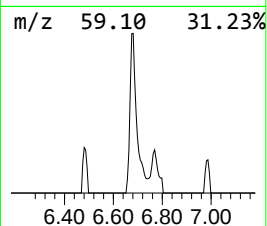
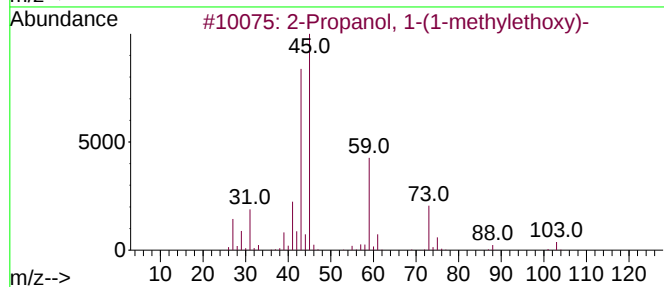
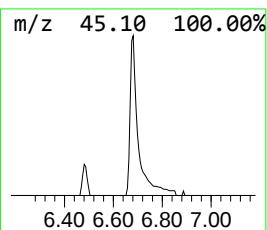
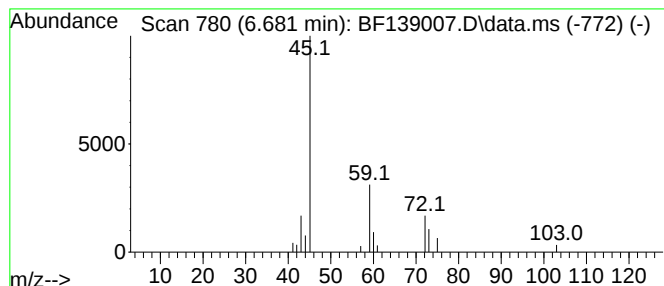
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 unknown6.681 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.681	2.39 ng	23647	1,4-Dichlorobenzene-d4	6.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(1-methylethoxy)-		118	C6H14O2	003944-36-3	38	
2	Ethyl ether		74	C4H10O	000060-29-7	38	
3	Ethane, 1,2-diethoxy-		118	C6H14O2	000629-14-1	9	
4	Diethyl carbitol		162	C8H18O3	000112-36-7	9	
5	2-Propanol, 1-propoxy-		118	C6H14O2	001569-01-3	9	



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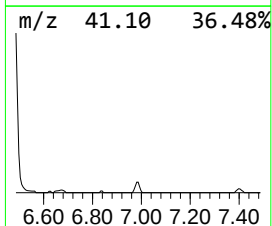
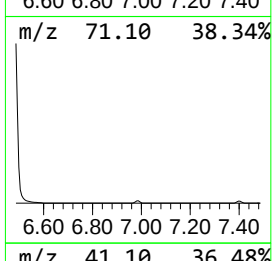
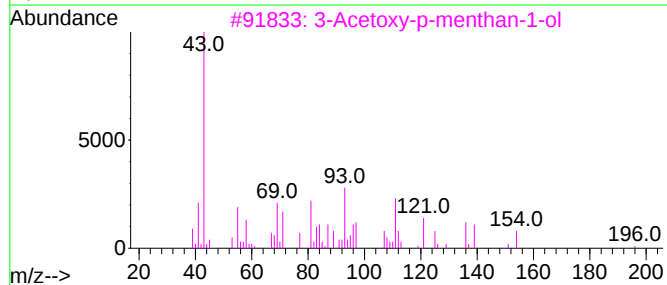
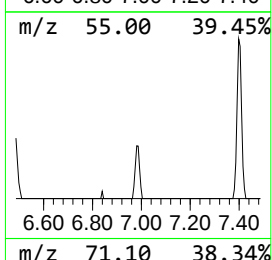
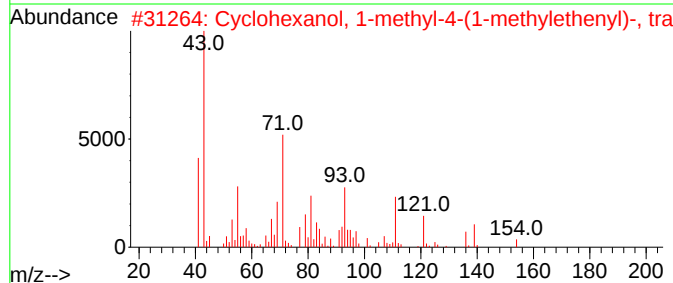
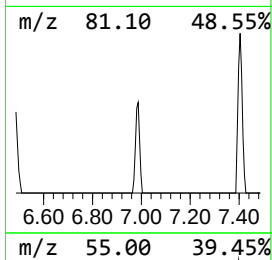
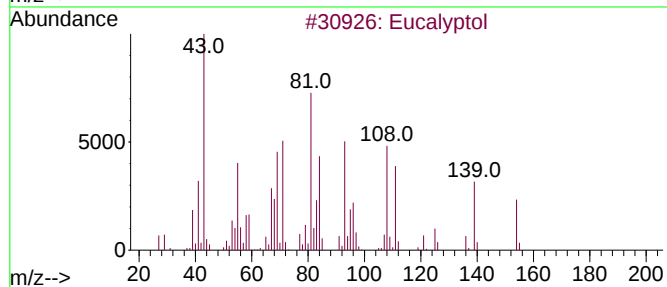
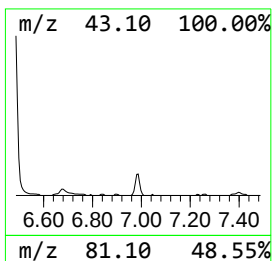
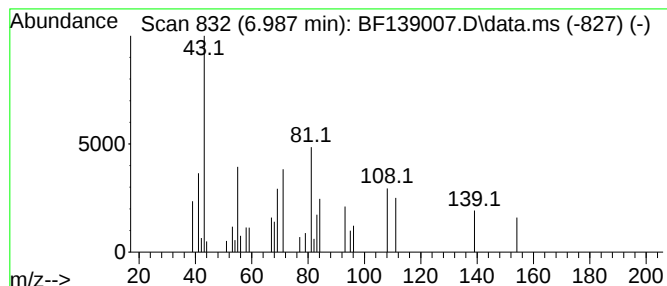
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Peak Number 4 Eucalyptol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.987	2.35 ng	23269	1,4-Dichlorobenzene-d4	6.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eucalyptol	154	C10H18O	000470-82-6	91
2			Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	007299-40-3	40
3			3-Acetoxy-p-menthan-1-ol	214	C12H22O3	058315-86-9	17
4			2-Isopropyl-5-methylhex-2-enal	154	C10H18O	035158-25-9	14
5			2-(3'-Oxo-butyl)-cyclopentanone	154	C9H14O2	1010143-37-2	14



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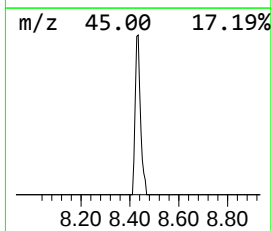
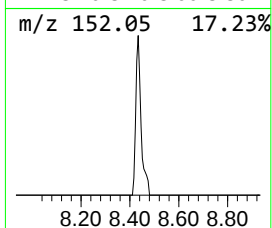
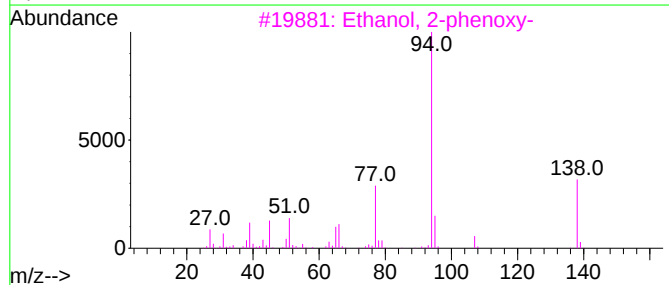
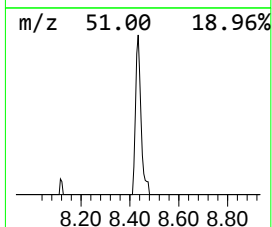
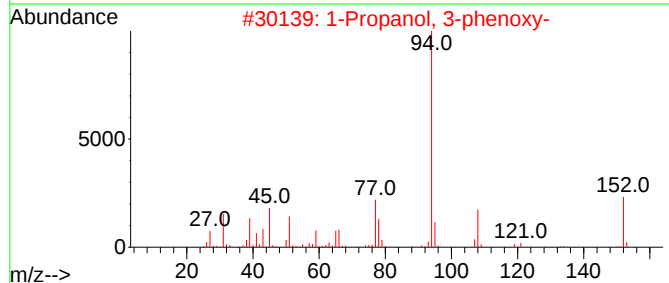
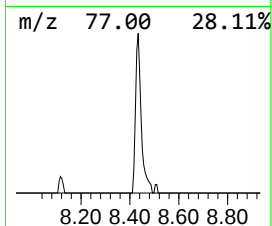
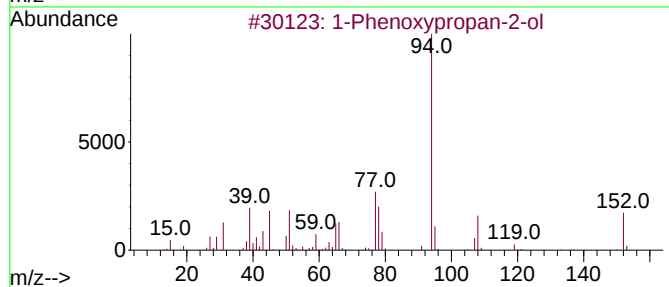
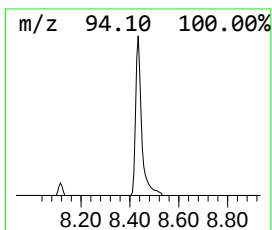
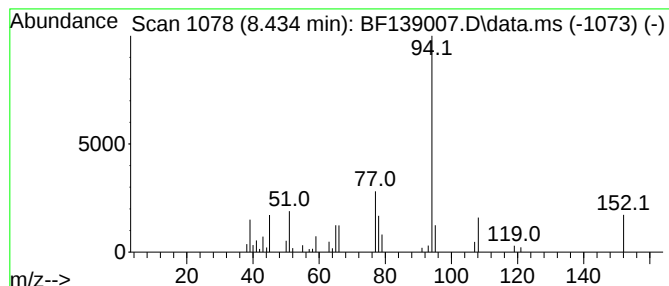
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.434	5.29 ng	55504	Naphthalene-d8	8.116

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	91
3		Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4		3-(.alpha.-Hydroxyethyl)-aniline	137	C8H11NO	002454-37-7	59
5		Benzene, (2-methylpropoxy)-	150	C10H14O	001126-75-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
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Instrument :
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ClientSampleId :
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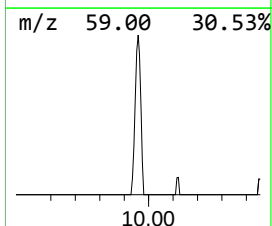
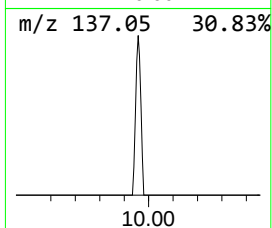
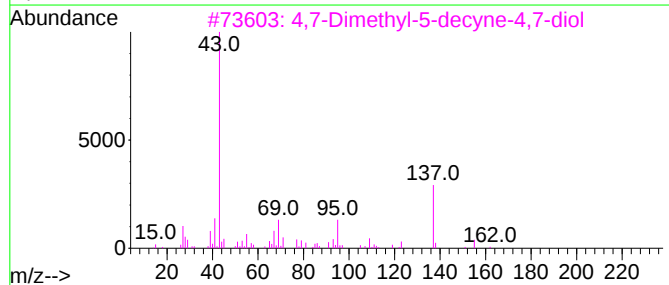
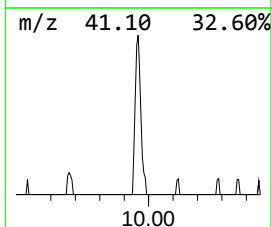
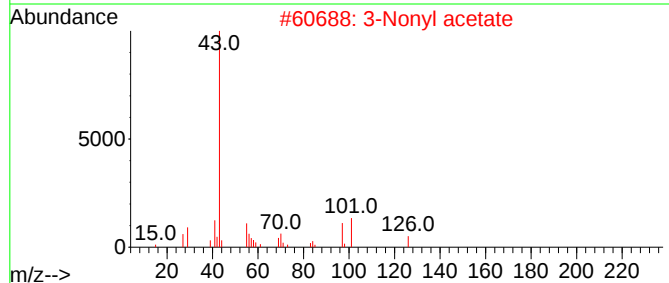
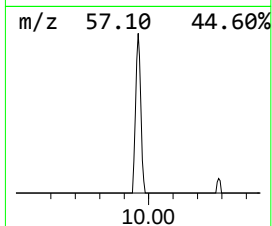
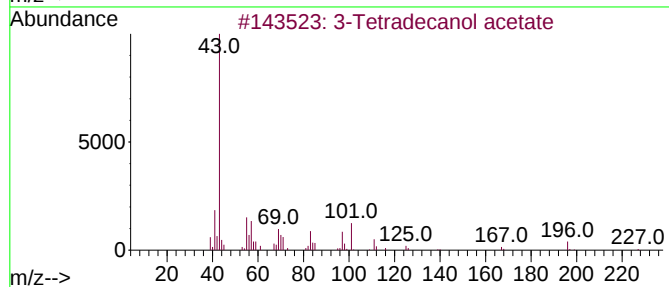
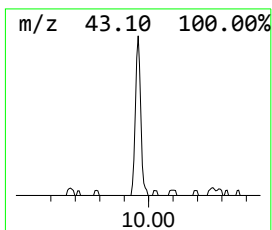
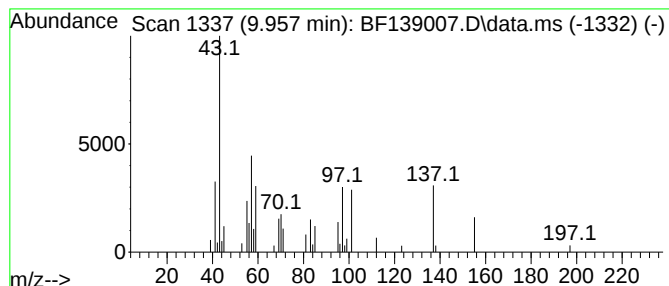
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown9.957 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.957	4.19 ng	40803	Acenaphthene-d10	9.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Tetradecanol acetate	256	C16H32O2	051354-23-5	16
2			3-Nonyl acetate	186	C11H22O2	1010429-16-2	12
3			4,7-Dimethyl-5-decyne-4,7-diol	198	C12H22O2	000126-87-4	10
4			7-Oxabicyclo[4.1.0]heptane, 1-me...	112	C7H12O	001713-33-3	10
5			2-Piperidinone, N-[4-bromo-n-but...	233	C9H16BrNO	195194-80-0	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
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Instrument :
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ClientSampleId :
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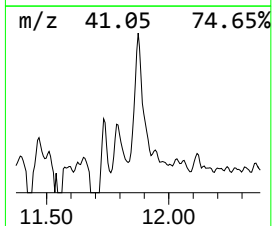
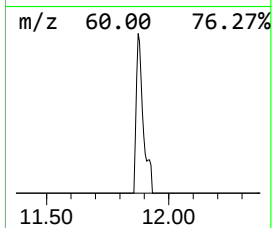
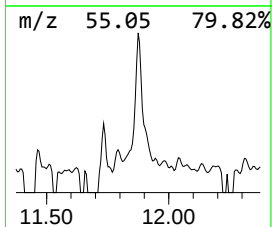
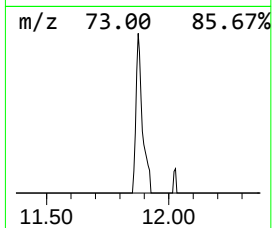
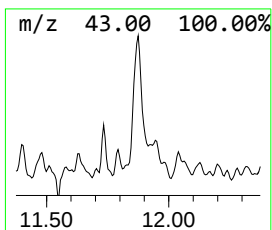
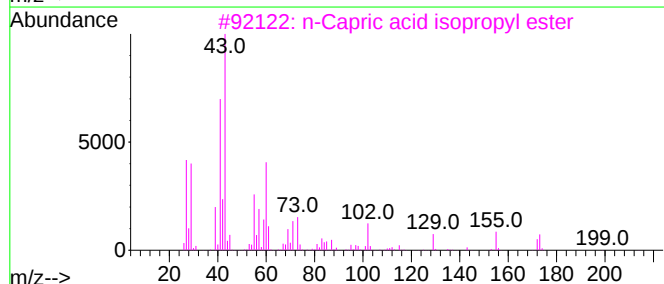
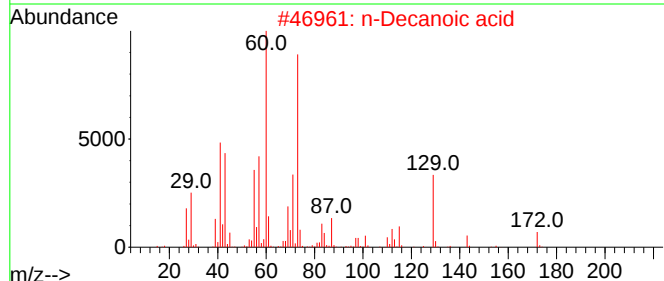
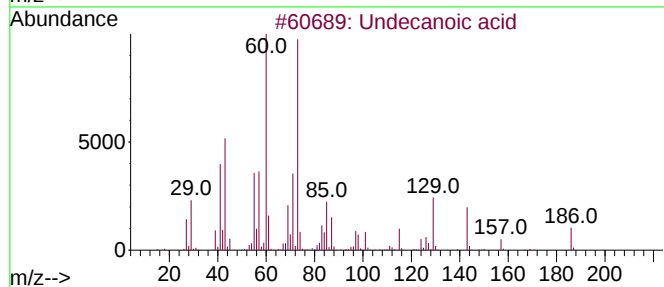
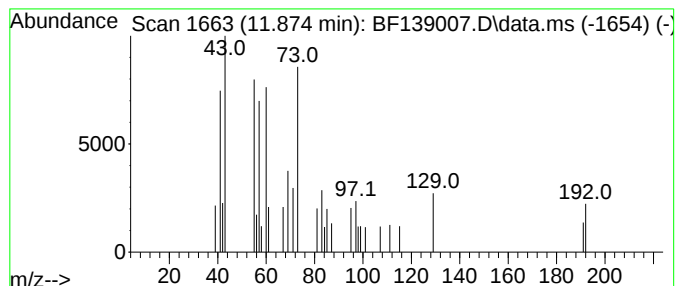
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Undecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.874	2.46 ng	29409	Phenanthrene-d10	11.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecanoic acid	186	C11H22O2	000112-37-8	52
2			n-Decanoic acid	172	C10H20O2	000334-48-5	43
3			n-Capric acid isopropyl ester	214	C13H26O2	002311-59-3	43
4			7-Methyloctanoic acid	158	C9H18O2	000693-19-6	38
5			Maltose	342	C12H22O11	000069-79-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

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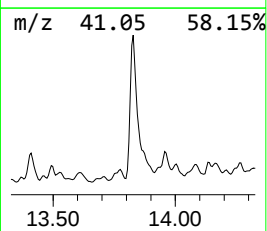
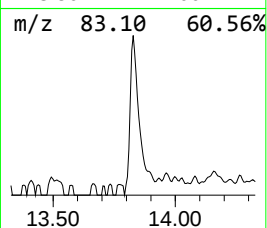
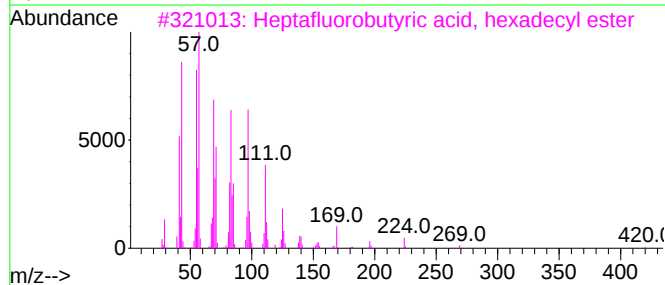
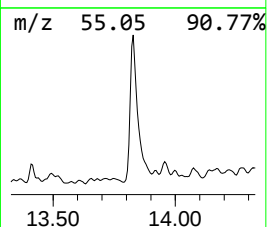
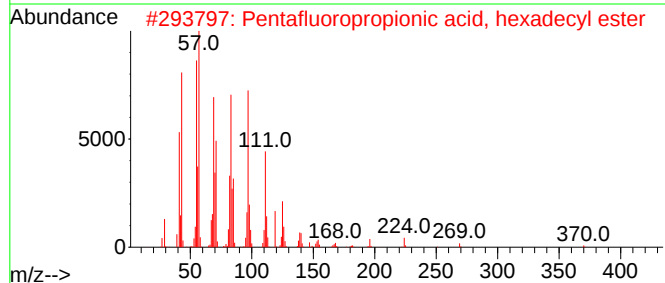
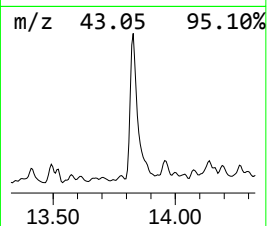
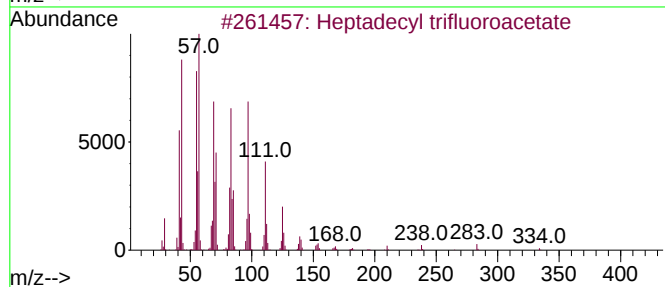
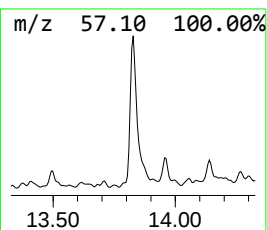
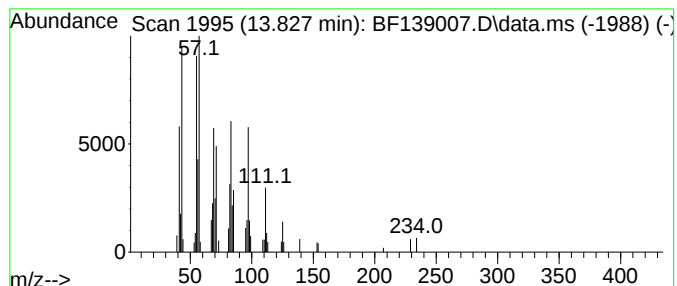
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecyl trifluoroacetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.827	8.99 ng	93390	Chrysene-d12	13.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91
2			Pentafluoropropionic acid, hexadecyl ester	388	C19H33F5O2	006222-07-7	91
3			Heptafluorobutyric acid, hexadecyl ester	438	C20H33F7O2	006385-15-5	91
4			Octacosanol	410	C28H58O	000557-61-9	91
5			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

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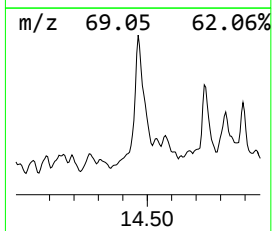
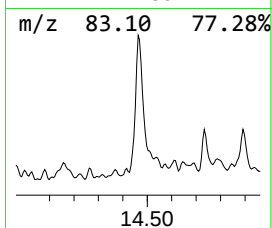
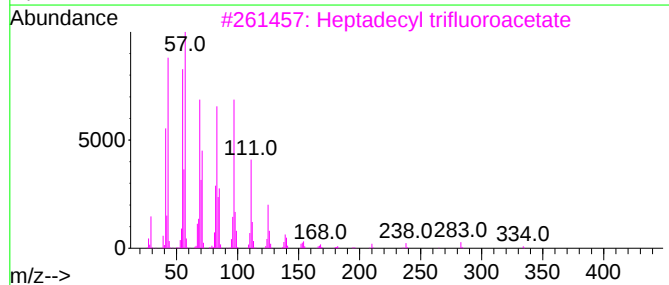
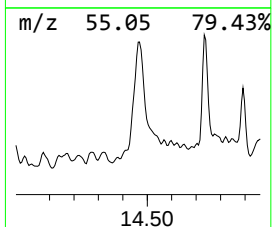
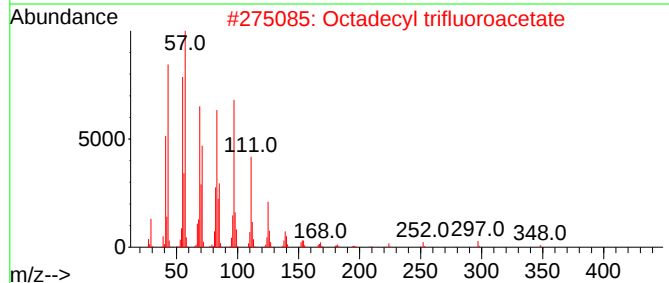
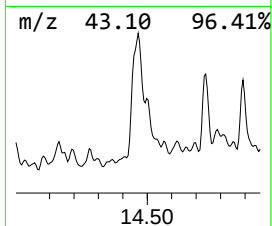
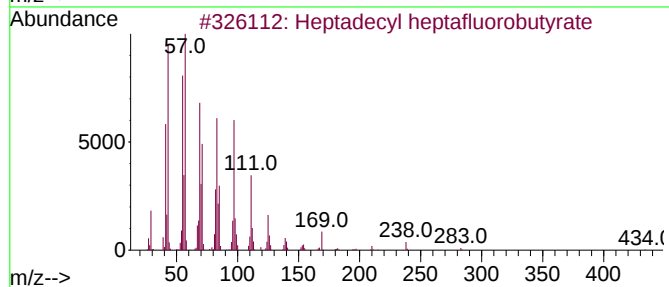
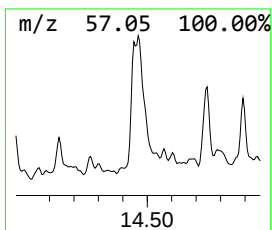
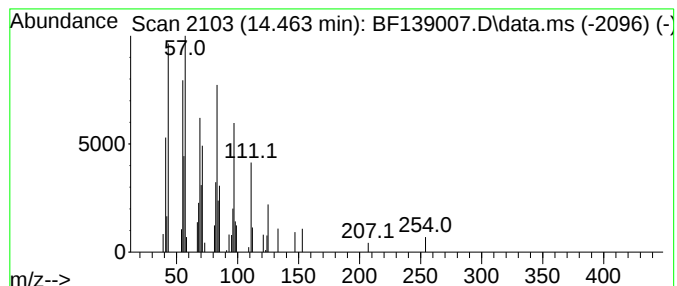
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecyl heptafluorobutyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	6.46 ng	67087	Chrysene-d12	13.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
2			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91
4			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
5			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
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Instrument :
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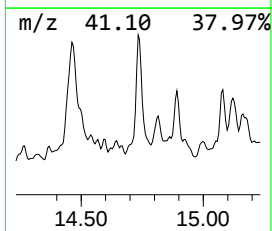
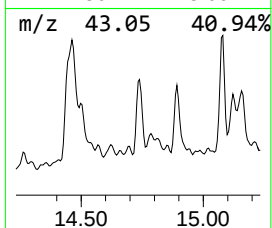
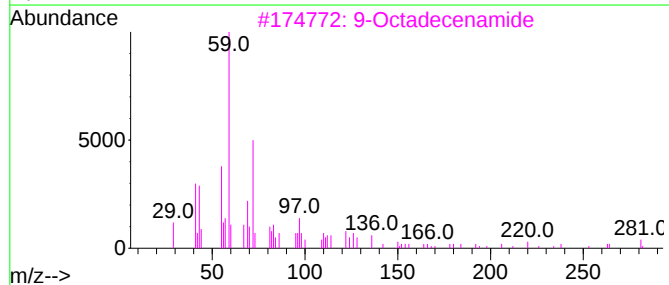
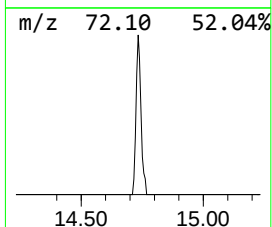
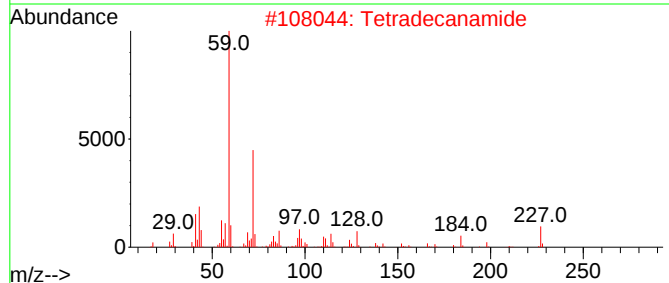
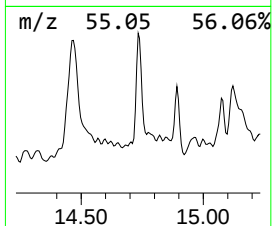
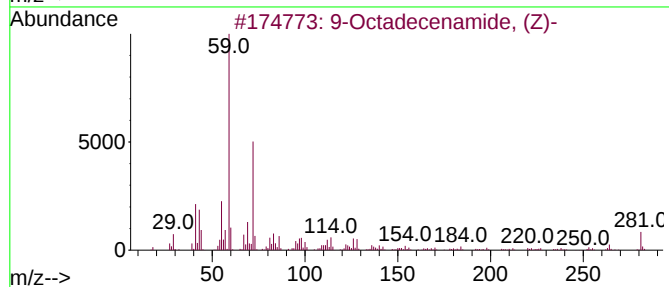
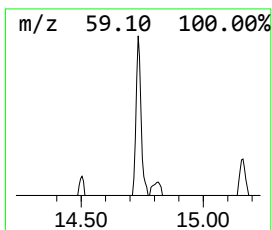
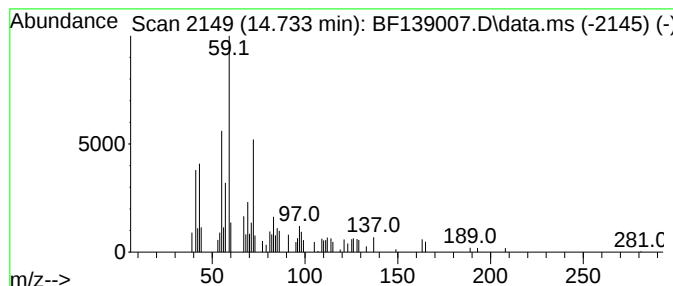
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.733	5.44 ng	34100	Perylene-d12	15.463

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
2		Tetradecanamide	227	C14H29NO	000638-58-4	80
3		9-Octadecenamide	281	C18H35NO	003322-62-1	64
4		Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1	53
5		Dodecanamide	199	C12H25NO	001120-16-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
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Sample : P3475-02
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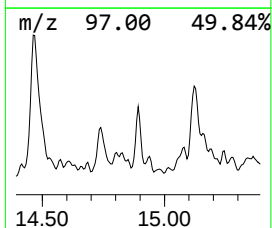
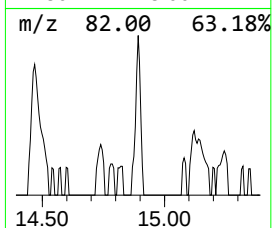
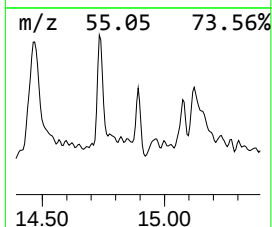
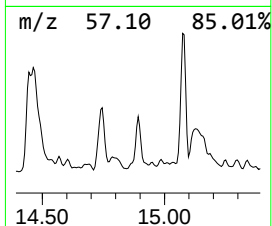
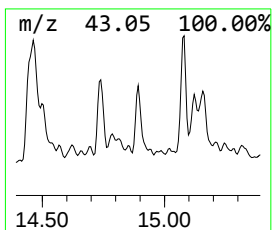
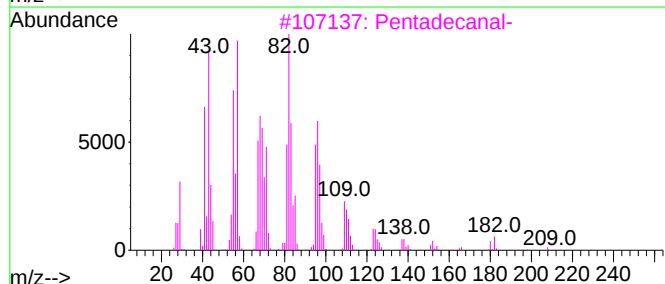
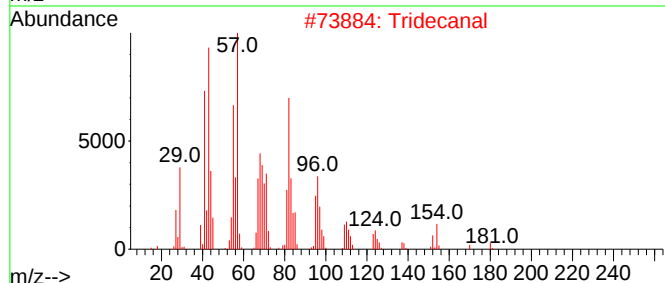
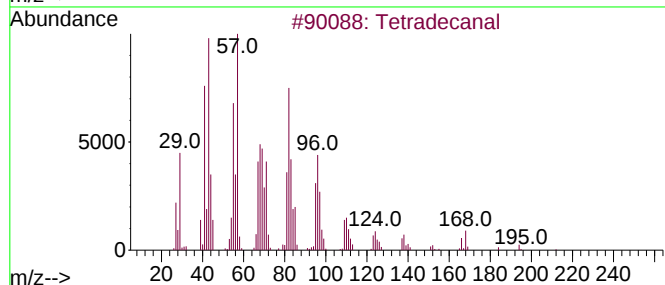
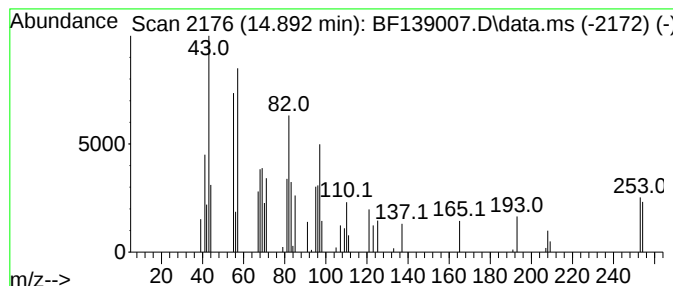
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetradecanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.892	2.88 ng	18078	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanal	212	C14H28O	000124-25-4	53
2			Tridecanal	198	C13H26O	010486-19-8	47
3			Pentadecanal-	226	C15H30O	002765-11-9	47
4			12-Octadecenal	266	C18H34O	056554-91-7	47
5			Octadecanal	268	C18H36O	000638-66-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
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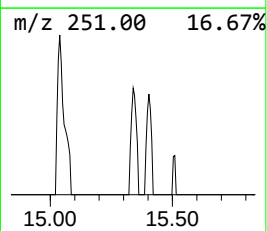
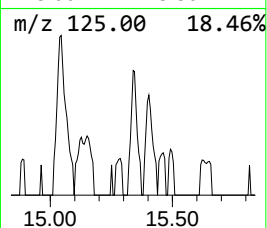
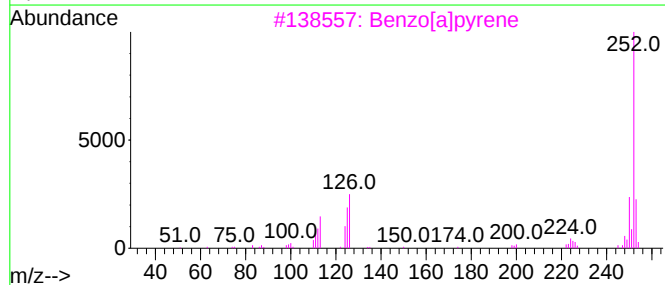
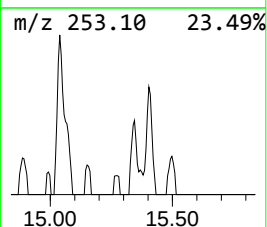
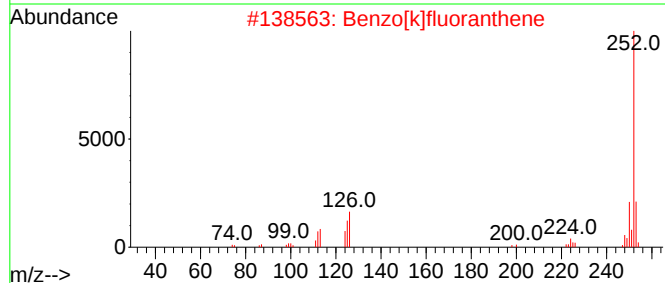
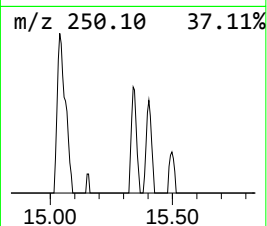
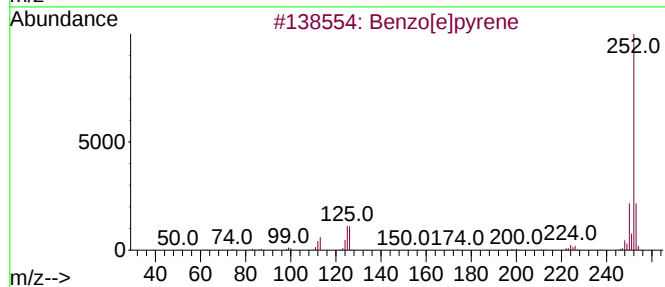
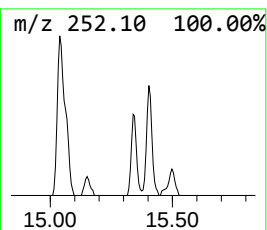
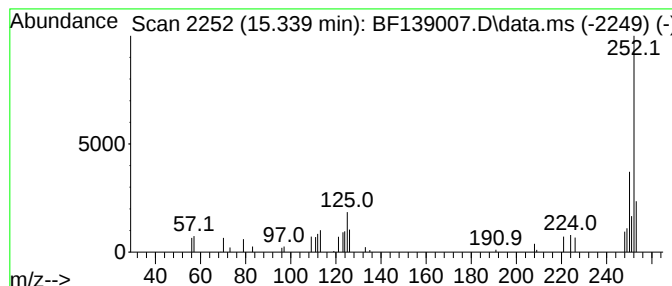
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.339	2.13 ng	13324	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	64
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	62
3			Benzo[a]pyrene	252	C20H12	000050-32-8	58
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	58
5			Perylene	252	C20H12	000198-55-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-866-N-S0-1.0-1.5-080524

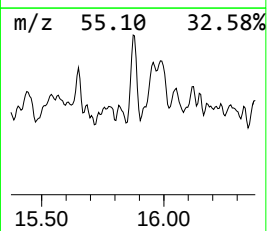
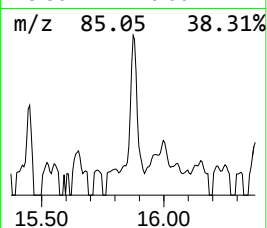
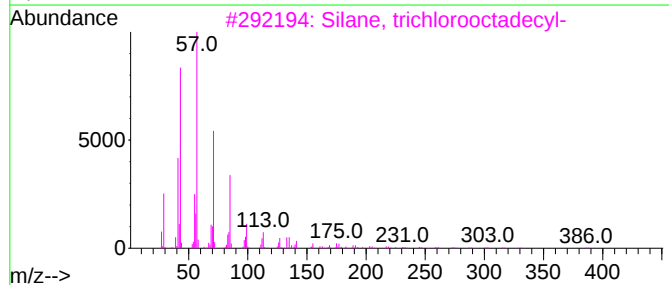
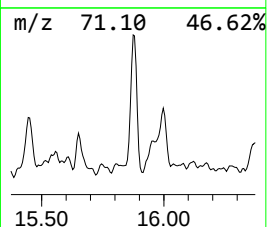
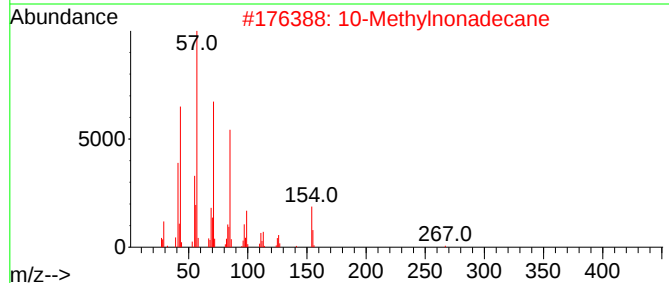
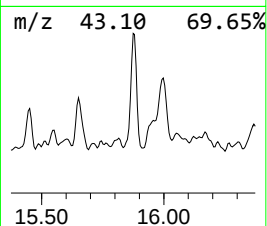
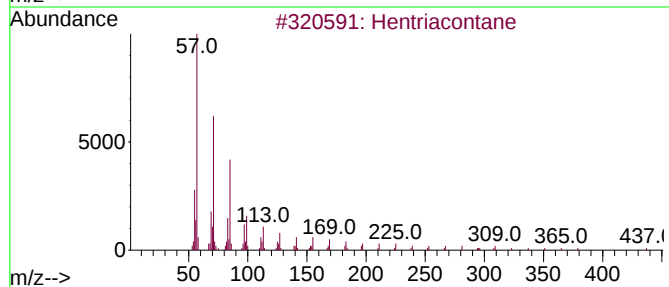
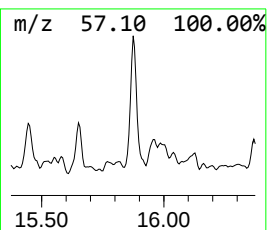
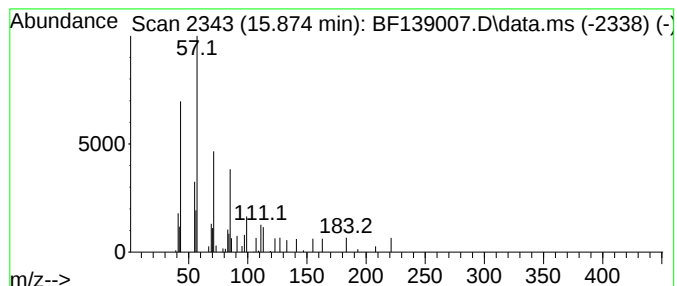
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hentriacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.874	3.34 ng	20949	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	80
2			10-Methylnonadecane	282	C20H42	056862-62-5	80
3			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	80
4			Hexacosane	366	C26H54	000630-01-3	80
5			Hexatriacontane	507	C36H74	000630-06-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-866-N-S0-1.0-1.5-080524

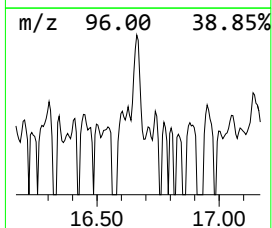
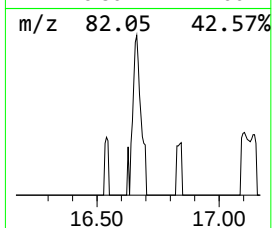
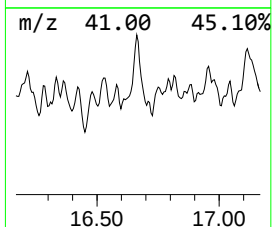
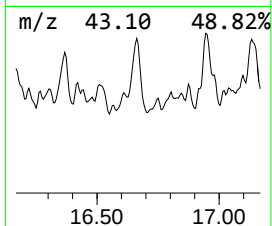
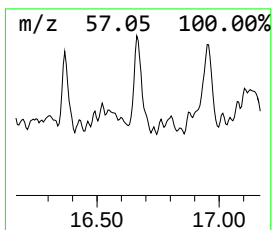
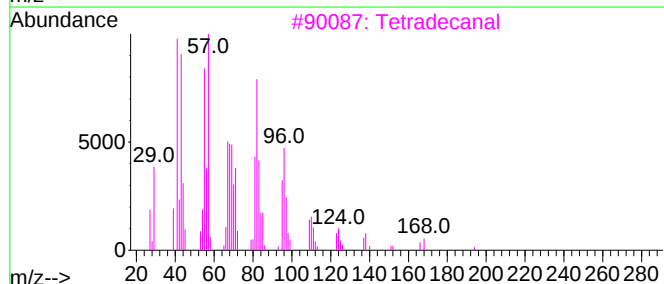
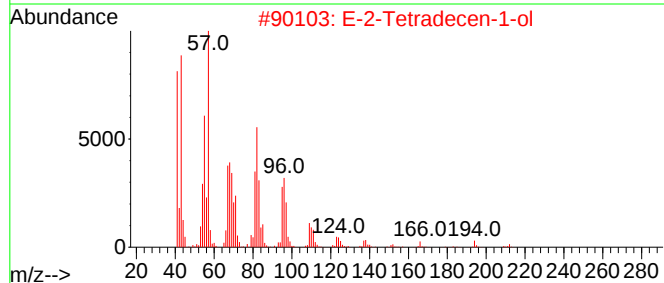
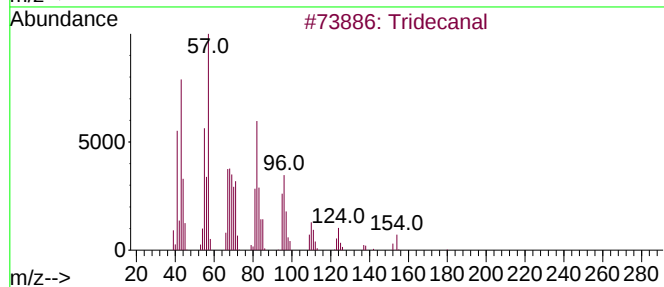
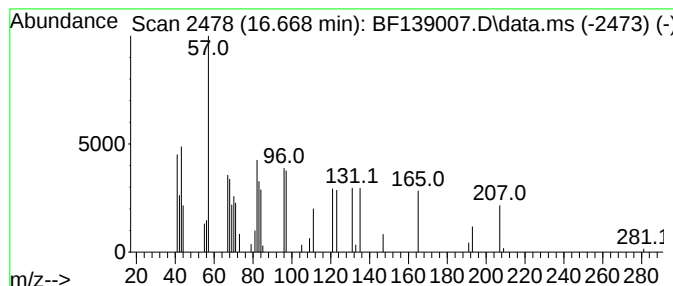
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown16.668 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.668	2.17 ng	13603	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanal	198	C13H26O	010486-19-8	22
2			E-2-Tetradecen-1-ol	212	C14H28O	1000130-83-7	14
3			Tetradecanal	212	C14H28O	000124-25-4	14
4			Z-2-Dodecenol	184	C12H24O	069064-36-4	10
5			Dodecanal	184	C12H24O	000112-54-9	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-866-N-S0-1.0-1.5-080524

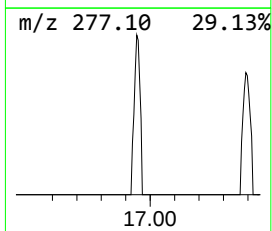
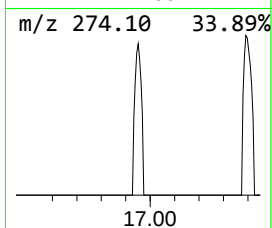
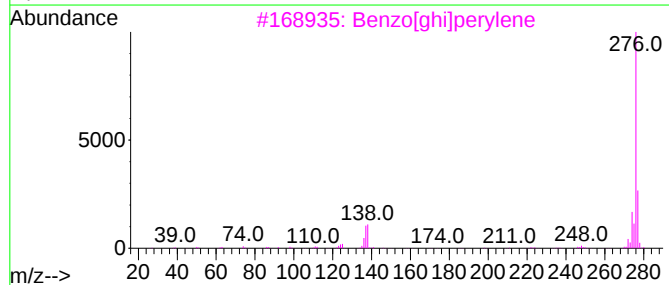
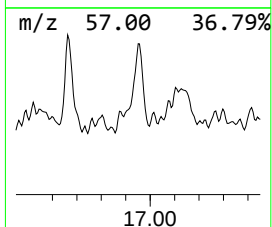
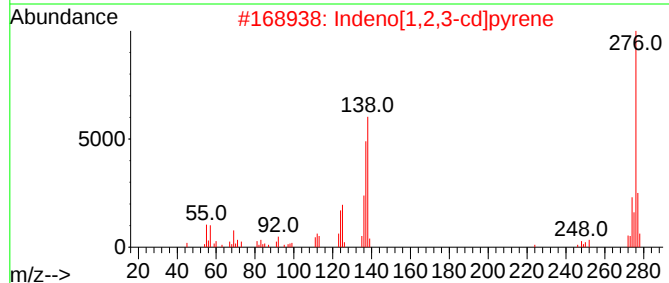
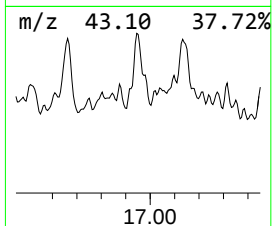
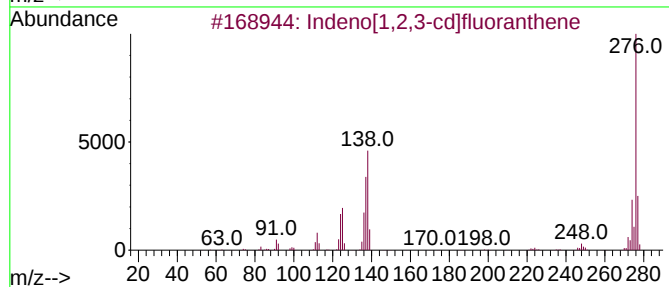
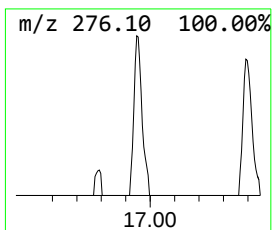
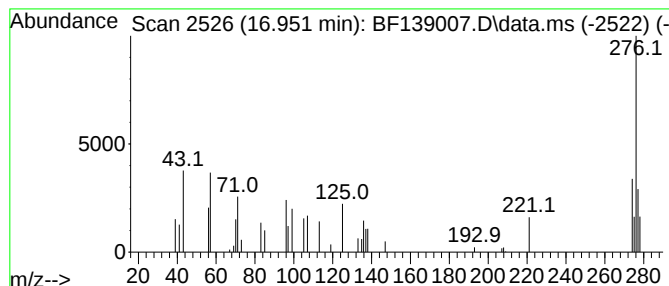
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown16.951 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.951	2.13 ng	13329	Perylene-d12	15.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	47
2			Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	47
3			Benzo[ghi]perylene	276	C22H12	000191-24-2	46
4			2-Amino-3-cyano-4,5-diphenylthio...	276	C17H12N2S	042160-34-9	43
5			Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF139007.D
Acq On : 14 Aug 2024 21:47
Operator : RC/JU
Sample : P3475-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-866-N-S0-1.0-1.5-080524

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
R-(-)-1,2-propa...	3.375	2.1	ng	20620	1	6.834	198133	20.0
2-Pentanone, 4-...	5.057	11.3	ng	111411	1	6.834	198133	20.0
unknown6.681	6.681	2.4	ng	23647	1	6.834	198133	20.0
Eucalyptol	6.987	2.4	ng	23269	1	6.834	198133	20.0
1-Phenoxypropan...	8.434	5.3	ng	55504	2	8.116	209677	20.0
unknown9.957	9.957	4.2	ng	40803	3	9.869	194605	20.0
Undecanoic acid	11.874	2.5	ng	29409	4	11.357	238722	20.0
Heptadecyl trif...	13.827	9.0	ng	93390	5	13.998	207679	20.0
Heptadecyl hept...	14.463	6.5	ng	67087	5	13.998	207679	20.0
9-Octadecenamid...	14.733	5.4	ng	34100	6	15.463	125400	20.0
Tetradecanal	14.892	2.9	ng	18078	6	15.463	125400	20.0
Benzo[e]pyrene	15.339	2.1	ng	13324	6	15.463	125400	20.0
Hentriacontane	15.874	3.3	ng	20949	6	15.463	125400	20.0
unknown16.668	16.668	2.2	ng	13603	6	15.463	125400	20.0
unknown16.951	16.951	2.1	ng	13329	6	15.463	125400	20.0