

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
 Data File : BF140641.D
 Acq On : 26 Nov 2024 14:55
 Operator : RC/JU
 Sample : P4860-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PH2-BOT-005

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-BF112124.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF140641.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.122	3	5	14	rVB	606370	703888	41.04%	5.408%
2	5.087	500	509	519	rBV	19315	32671	1.90%	0.251%
3	5.498	573	579	594	rBV	965519	1282769	74.80%	9.856%
4	6.504	745	750	774	rBV	1030272	1372157	80.01%	10.543%
5	6.869	807	812	819	rBV	356677	430326	25.09%	3.306%
6	7.434	902	908	919	rBV	648844	887376	51.74%	6.818%
7	7.686	946	951	963	rBV2	10719	22096	1.29%	0.170%
8	8.151	1025	1030	1045	rBV	410981	535824	31.24%	4.117%
9	9.222	1207	1212	1223	rBV	1353393	1715039	100.00%	13.178%
10	9.904	1323	1328	1345	rVB	454879	590370	34.42%	4.536%
11	10.616	1444	1449	1458	rBV	217506	284390	16.58%	2.185%
12	10.698	1458	1463	1478	rVV	624413	837670	48.84%	6.436%
13	10.927	1497	1502	1512	rVB	34059	46679	2.72%	0.359%
14	11.398	1576	1582	1590	rBV	459145	602196	35.11%	4.627%
15	11.916	1666	1670	1684	rBV	82658	147459	8.60%	1.133%
16	12.416	1751	1755	1761	rBV2	15298	22741	1.33%	0.175%
17	12.645	1786	1794	1799	rBV	47372	72175	4.21%	0.555%
18	12.980	1845	1851	1861	rBV	1284783	1631156	95.11%	12.533%
19	13.204	1884	1889	1893	rBV	14086	20043	1.17%	0.154%
20	13.545	1943	1947	1953	rBV	24588	33211	1.94%	0.255%
21	13.874	1998	2003	2018	rBV	82803	161630	9.42%	1.242%
22	14.010	2022	2026	2028	rVV	25533	34821	2.03%	0.268%
23	14.045	2028	2032	2045	rVB	282281	389191	22.69%	2.990%
24	14.192	2053	2057	2062	rVB	35892	45459	2.65%	0.349%
25	14.504	2106	2110	2123	rBV	48562	80254	4.68%	0.617%
26	14.815	2158	2163	2169	rVB	34523	48570	2.83%	0.373%
27	15.157	2216	2221	2226	rVB	47975	65806	3.84%	0.506%
28	15.539	2280	2286	2301	rVB	268191	443348	25.85%	3.406%
29	15.980	2357	2361	2368	rVB	23818	39891	2.33%	0.307%
30	16.245	2399	2406	2423	rBV7	43904	147265	8.59%	1.132%
31	16.445	2433	2440	2445	rBV2	73348	142864	8.33%	1.098%
32	16.498	2446	2449	2464	rVB4	19590	49181	2.87%	0.378%
33	17.439	2605	2609	2620	rVB9	21092	50352	2.94%	0.387%
34	17.651	2641	2645	2658	rVB9	17542	45973	2.68%	0.353%

Sum of corrected areas: 13014841

Instrument :

BNA_F

ClientSampleId :

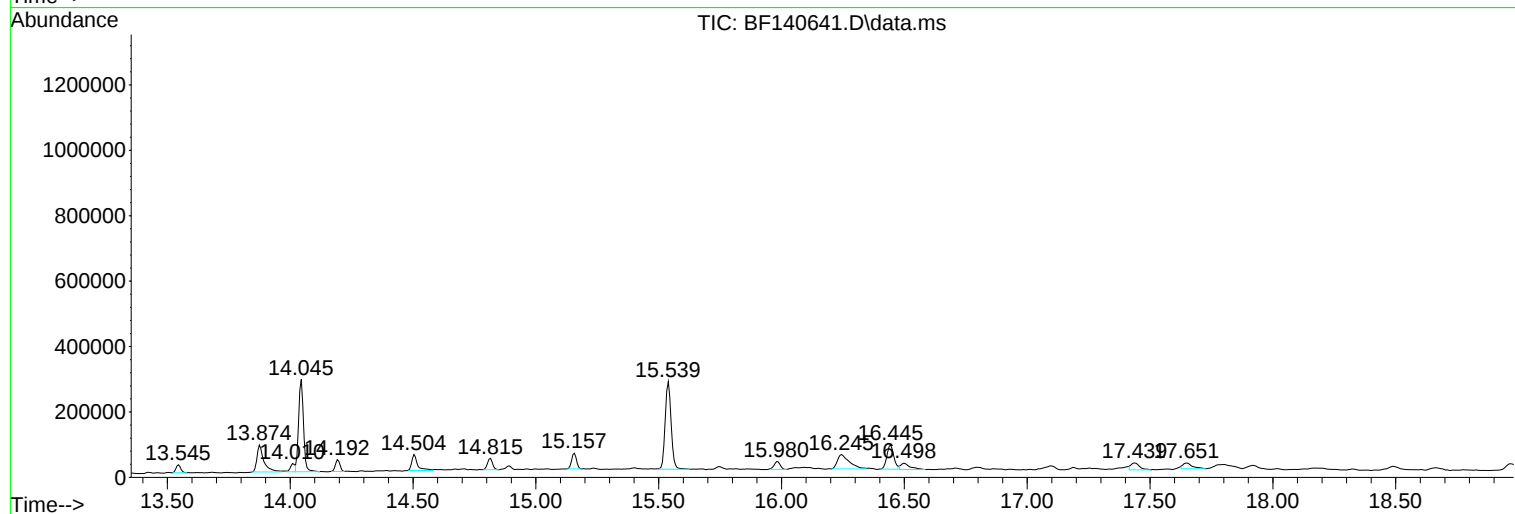
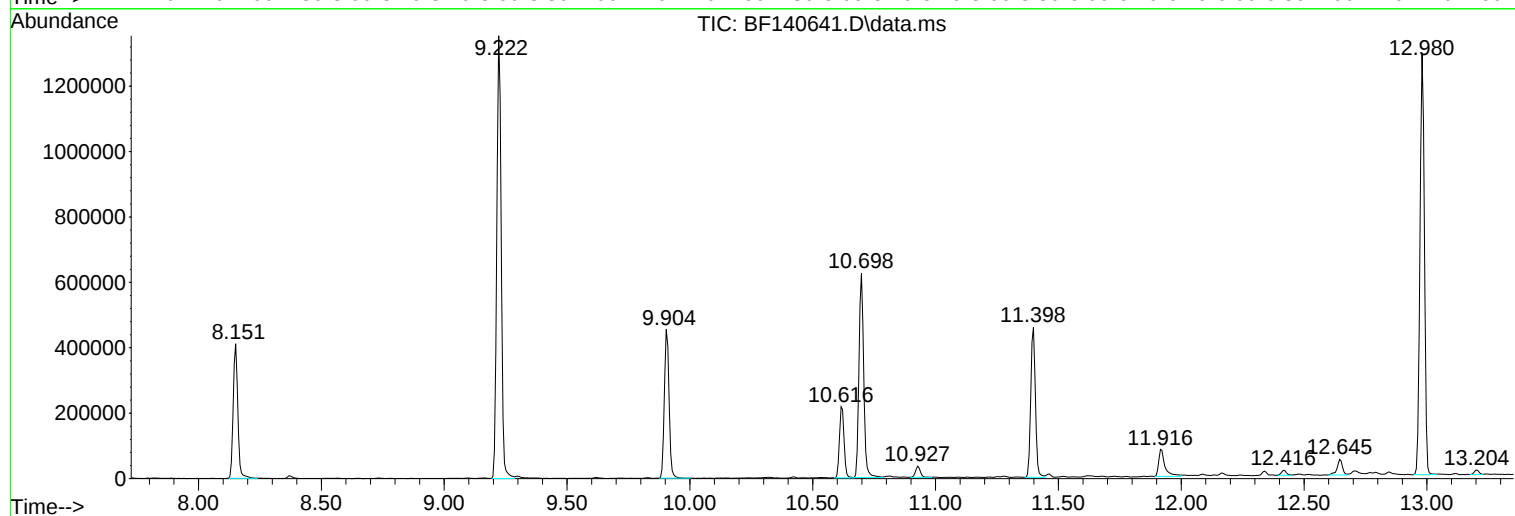
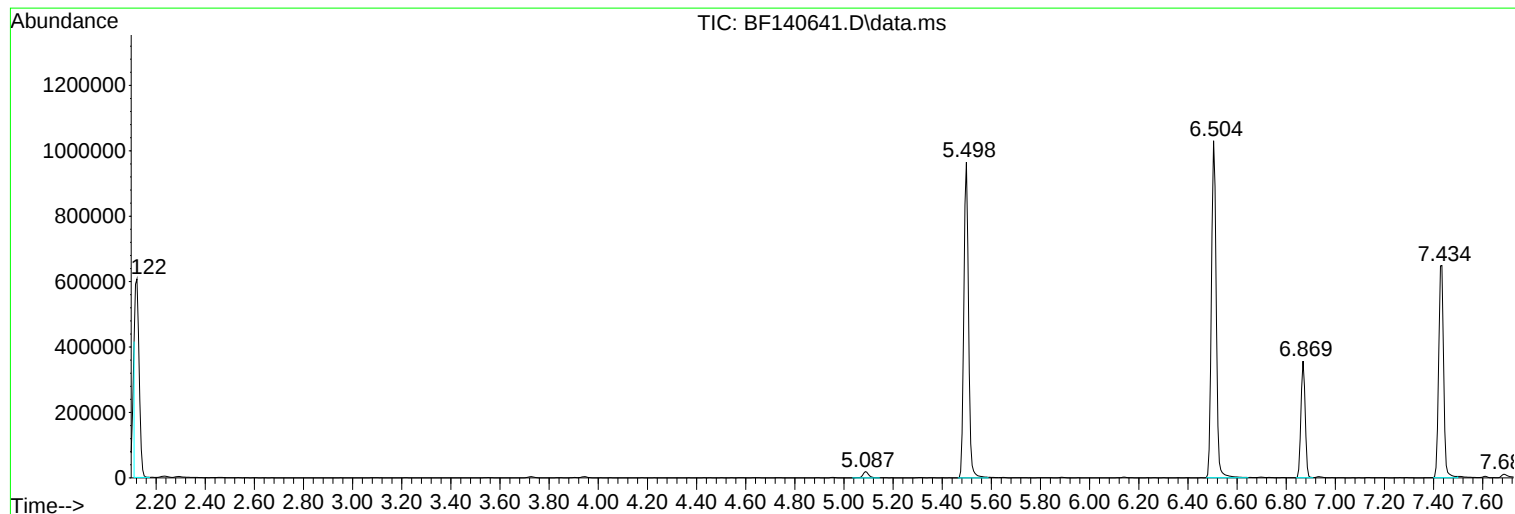
PH2-BOT-005

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.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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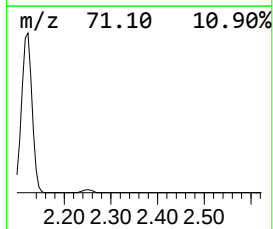
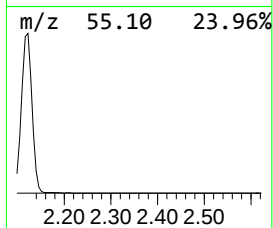
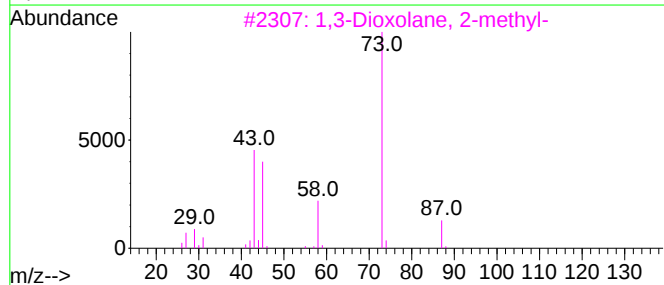
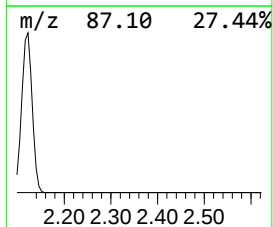
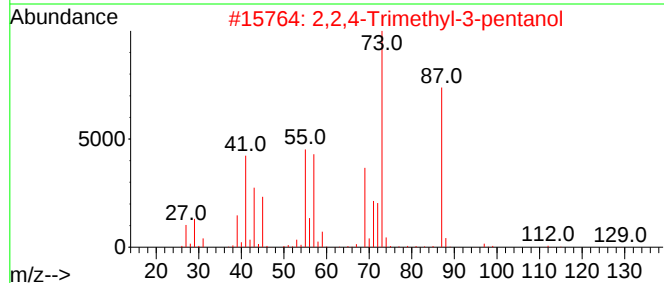
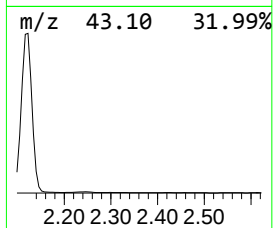
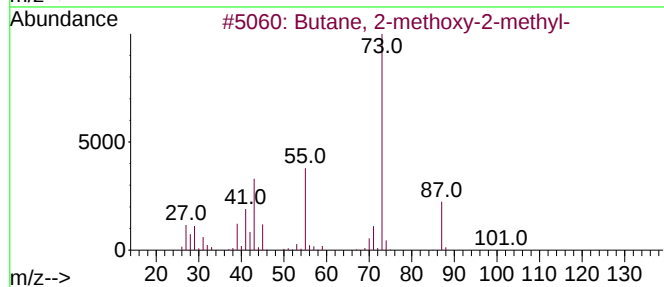
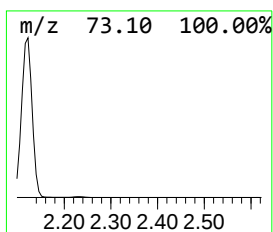
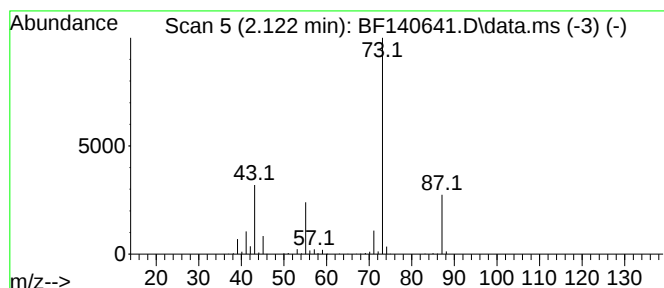
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.122	32.71 ng	703888	1,4-Dichlorobenzene-d4	6.869

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	25
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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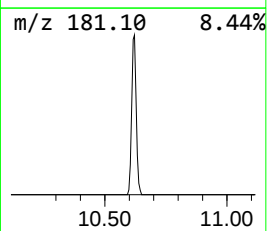
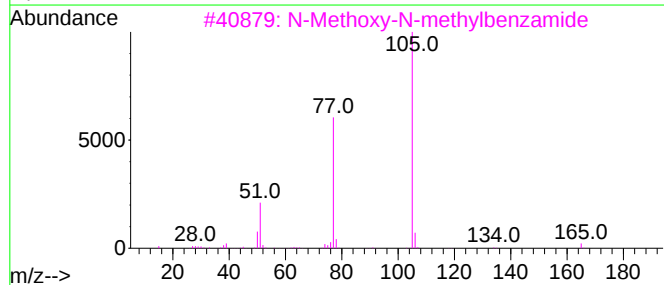
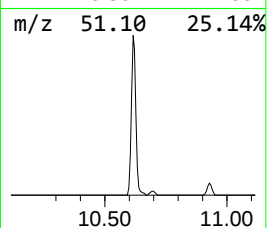
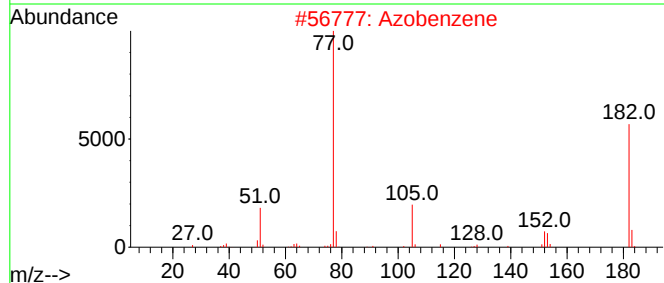
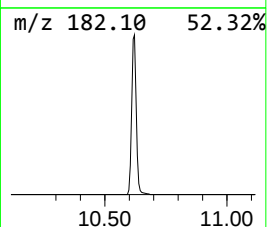
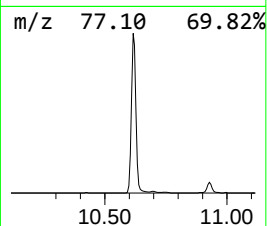
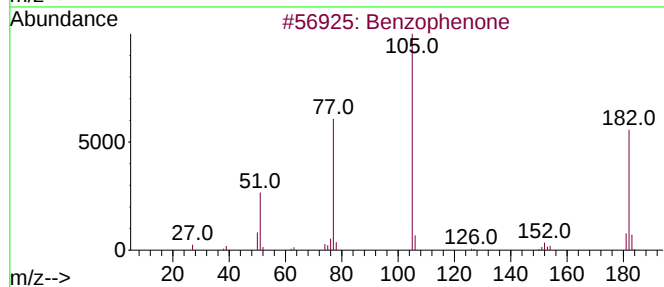
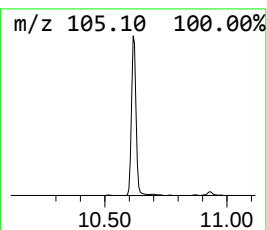
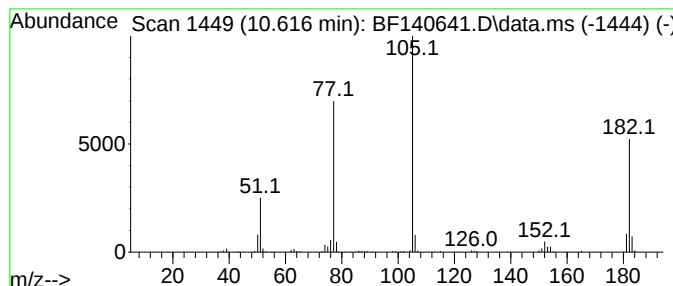
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.616	9.63 ng	284390	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Azobenzene	182	C12H10N2	000103-33-3	59
3			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	49
4			N-Methylbenzamide, N-trifluoroac...	231	C10H8F3NO2	1000446-91-5	47
5			Vinyl benzoate	148	C9H8O2	000769-78-8	47



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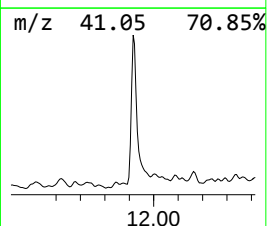
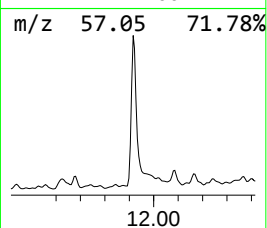
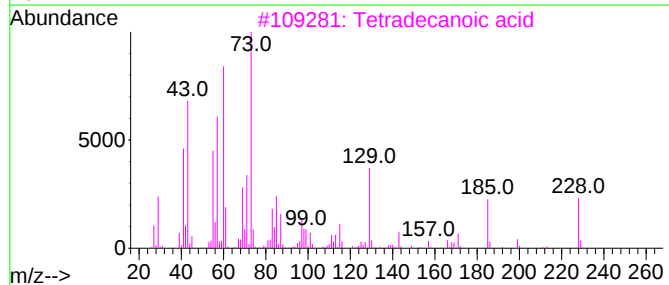
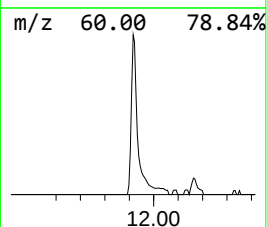
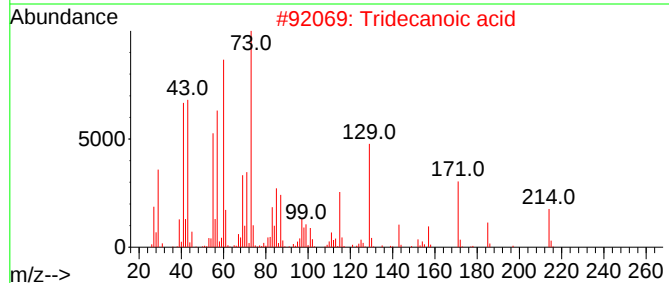
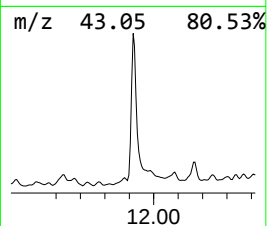
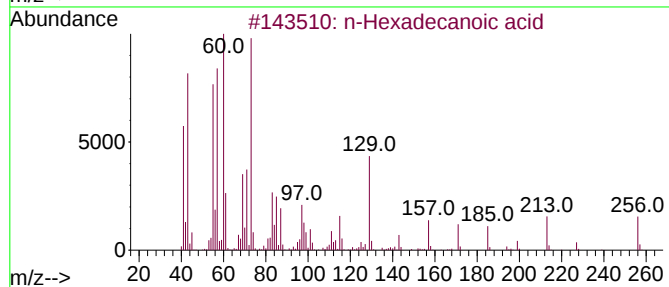
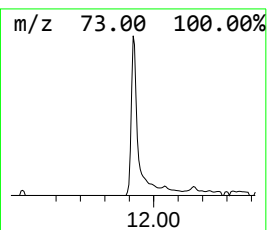
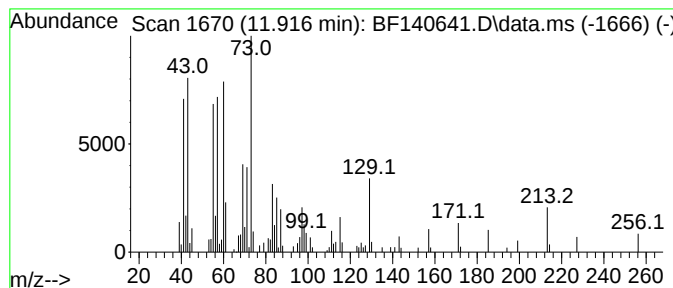
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	4.90 ng	147459	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			Tridecanoic acid	214	C13H26O2	000638-53-9	95
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	74
5			Oxalic acid, dodecyl propyl ester	300	C17H32O4	1000309-26-5	25



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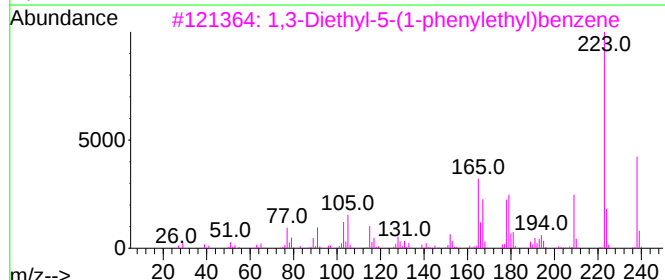
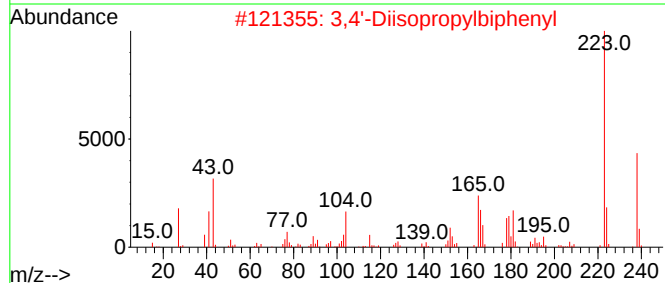
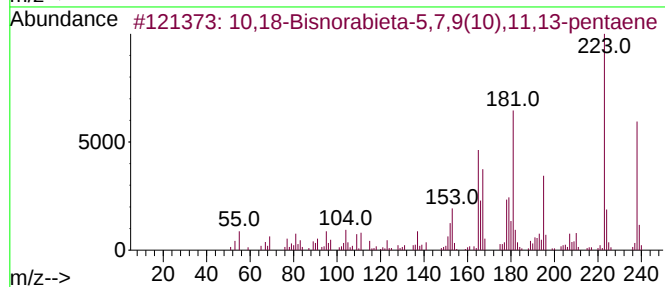
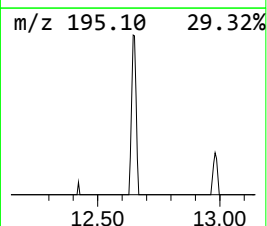
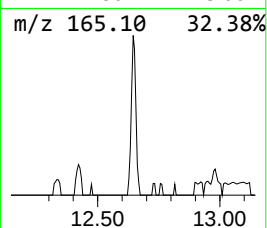
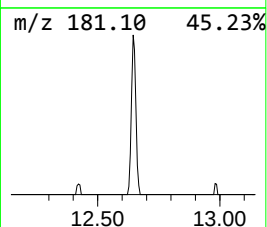
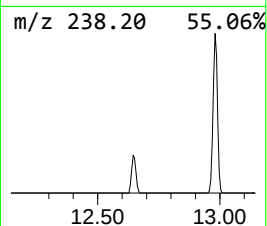
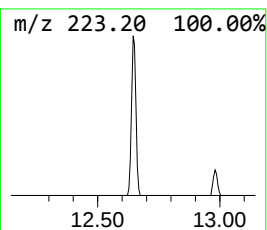
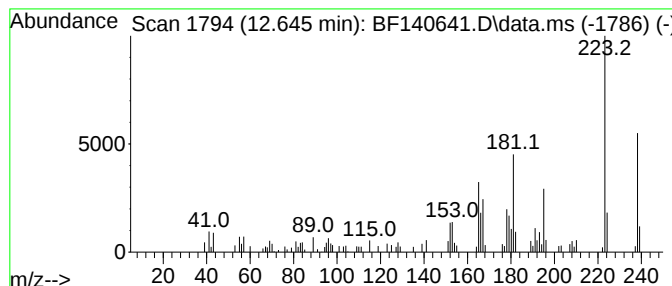
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.645	2.40 ng	72175	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99		
2	3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	89		
3	1,3-Diethyl-5-(1-phenylethyl)ben...	238	C18H22	1000482-32-6	72		
4	3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	49		
5	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	49		



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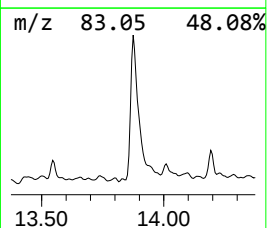
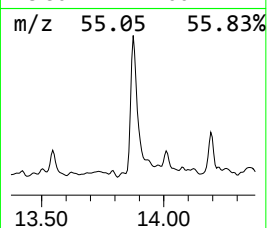
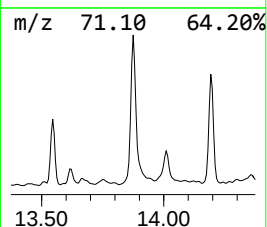
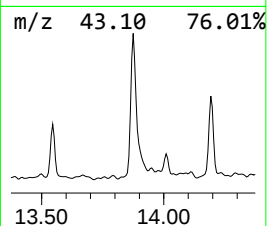
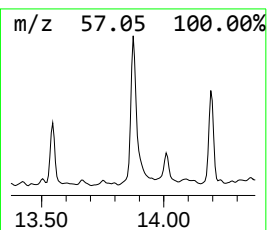
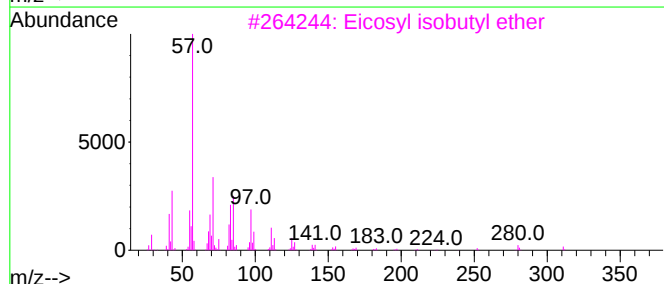
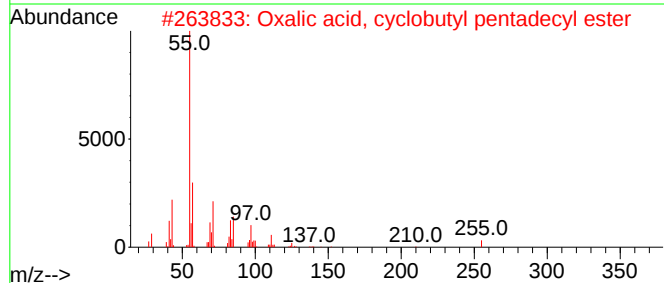
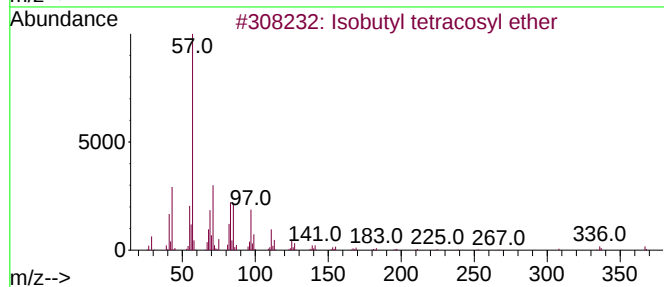
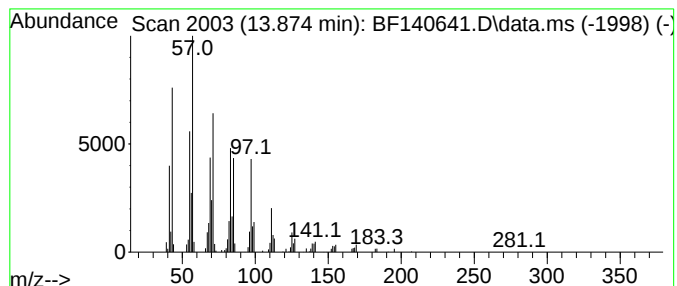
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Isobutyl tetracosyl ether Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	8.31 ng	161630	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl tetracosyl ether	410	C28H58O	1000406-33-2	91
2			Oxalic acid, cyclobutyl pentadec...	354	C21H38O4	1010309-70-5	90
3			Eicosyl isobutyl ether	354	C24H50O	1000406-33-0	90
4			1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	83
5			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	74



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Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
PH2-BOT-005

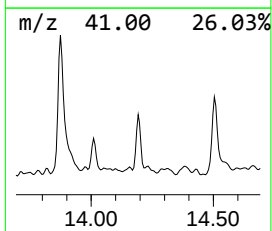
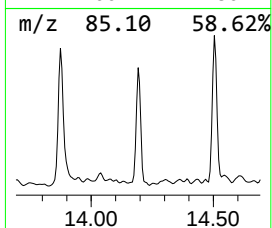
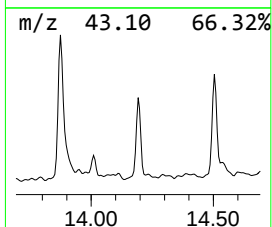
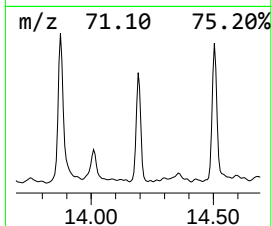
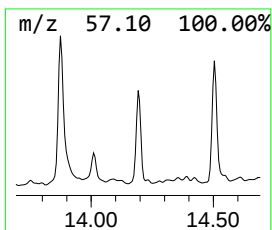
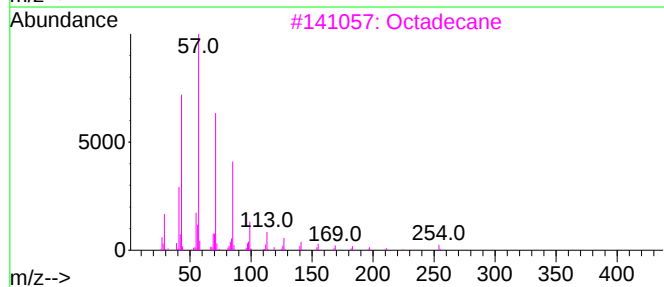
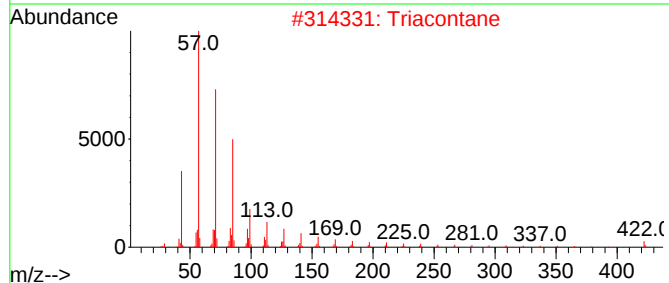
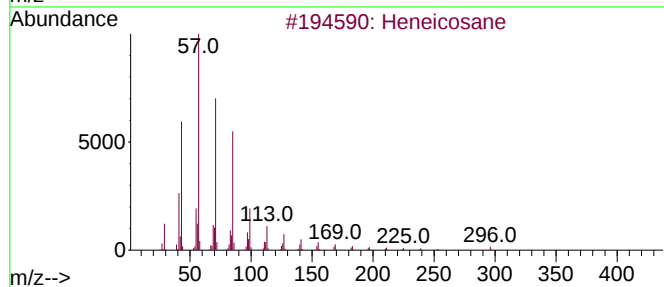
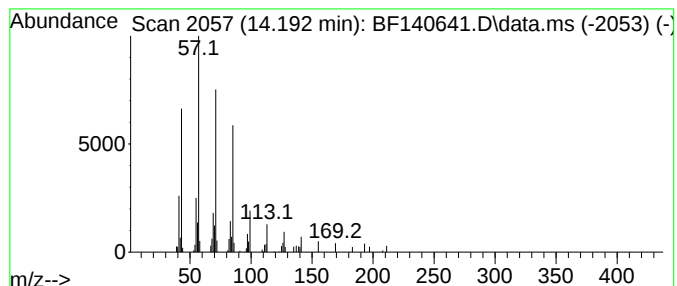
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.192	2.34 ng	45459	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Triacontane	422	C30H62	000638-68-6	91
3			Octadecane	254	C18H38	000593-45-3	91
4			Octacosane	394	C28H58	000630-02-4	90
5			Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1010115-66-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
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Acq On : 26 Nov 2024 14:55
Operator : RC/JU
Sample : P4860-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

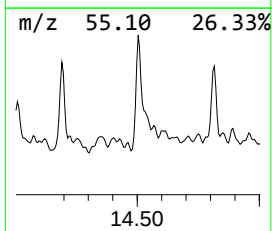
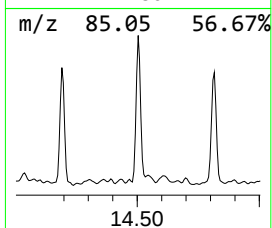
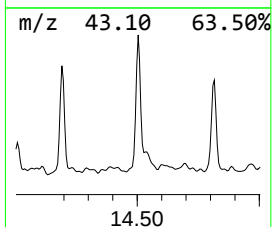
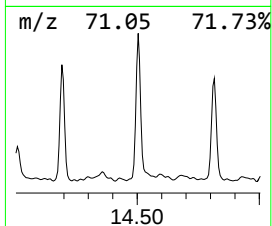
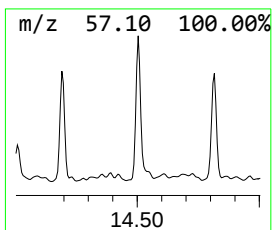
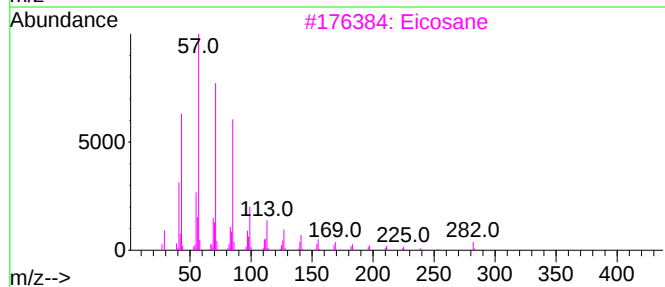
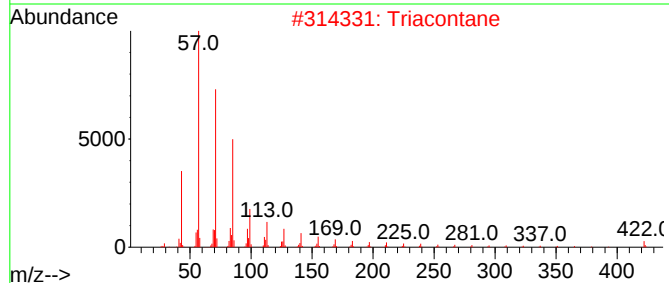
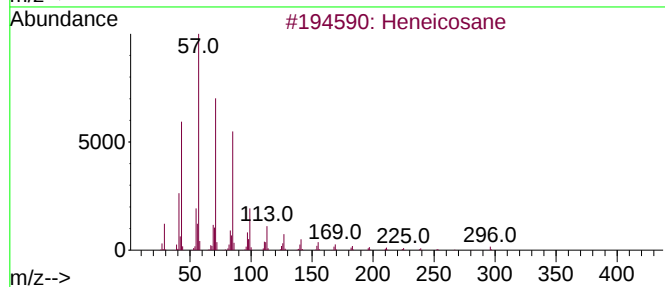
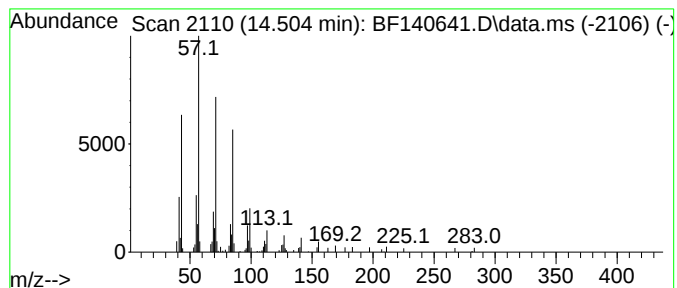
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Triacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.504	4.12 ng	80254	Chrysene-d12	14.045	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Triacontane	422	C30H62	000638-68-6	91
3	Eicosane	282	C20H42	000112-95-8	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
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Sample : P4860-10
Misc :
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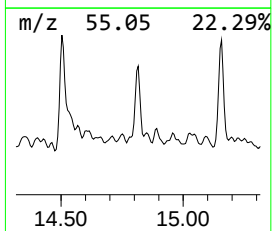
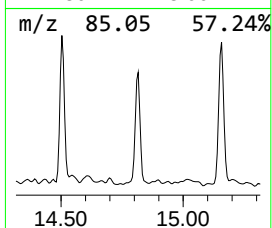
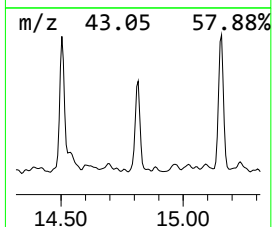
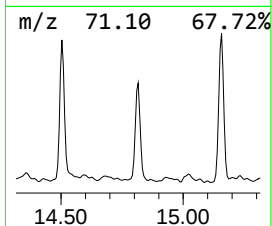
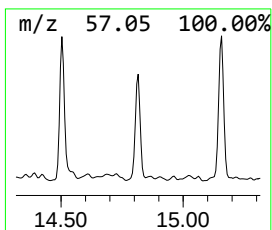
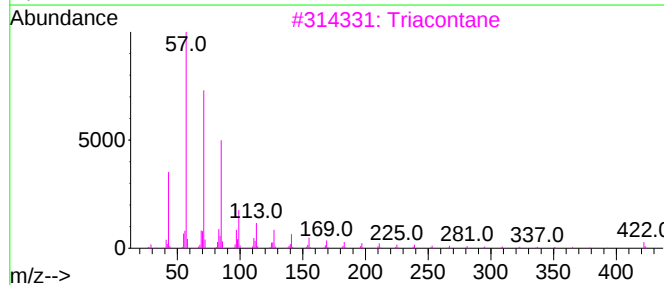
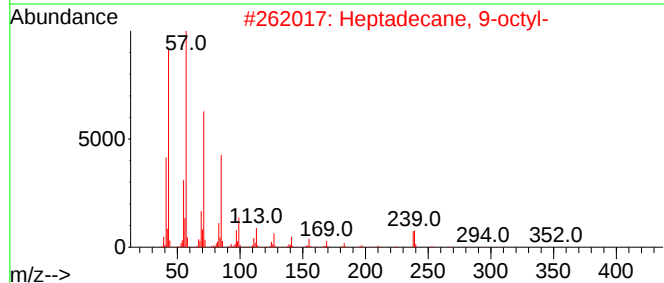
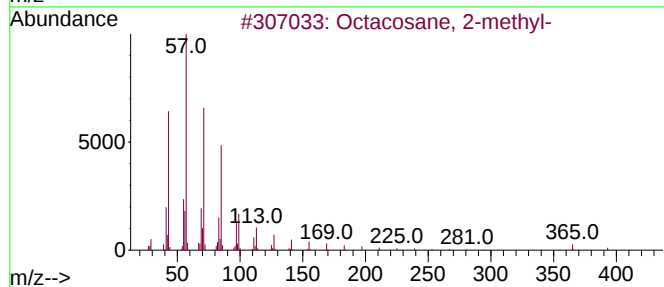
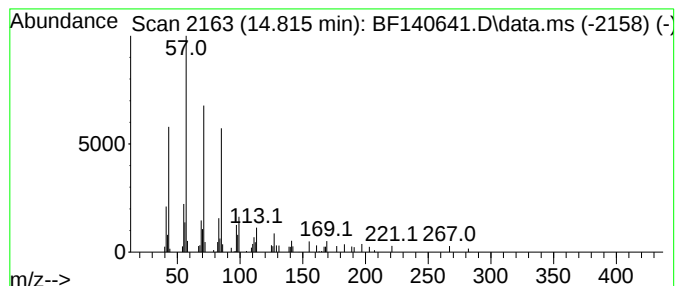
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.815	2.19 ng	48570	Perylene-d12	15.539	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Triacontane	422	C30H62	000638-68-6	91
4	Octacosane	394	C28H58	000630-02-4	91
5	Octadecane	254	C18H38	000593-45-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
Data File : BF140641.D
Acq On : 26 Nov 2024 14:55
Operator : RC/JU
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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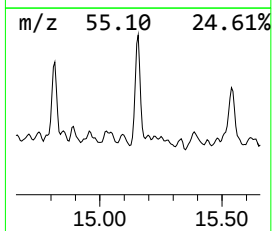
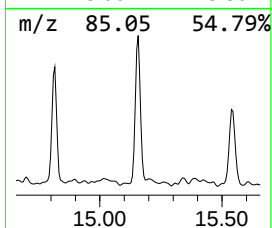
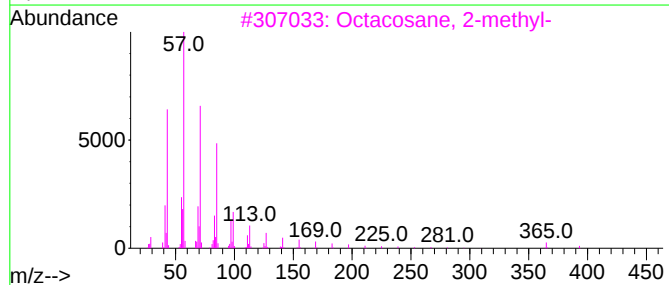
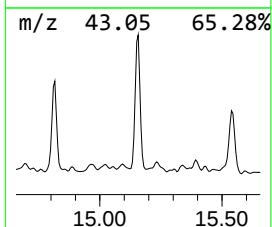
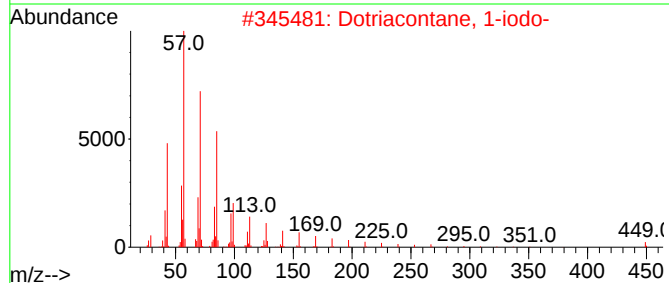
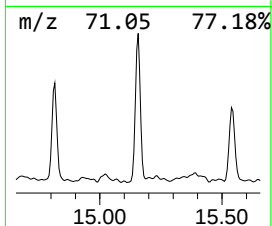
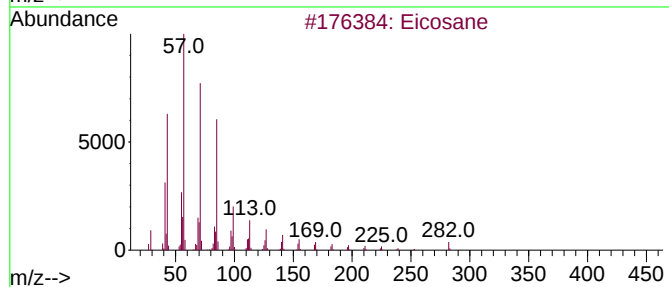
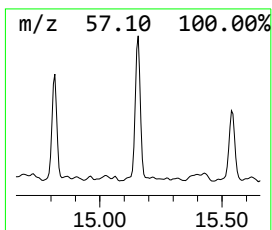
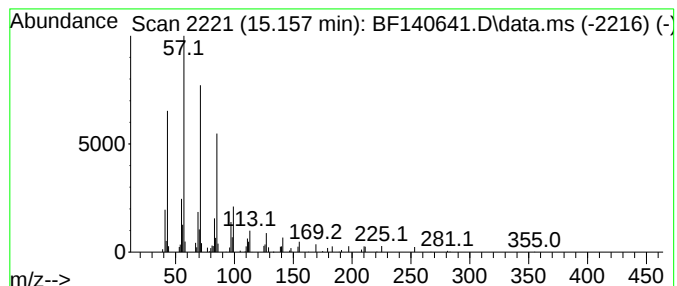
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.157	2.97 ng	65806	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
Data File : BF140641.D
Acq On : 26 Nov 2024 14:55
Operator : RC/JU
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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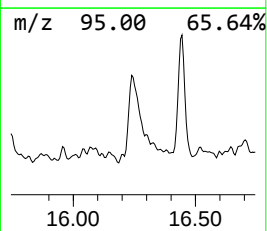
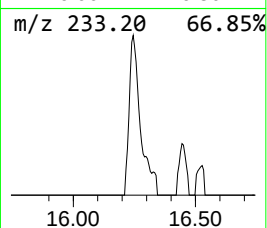
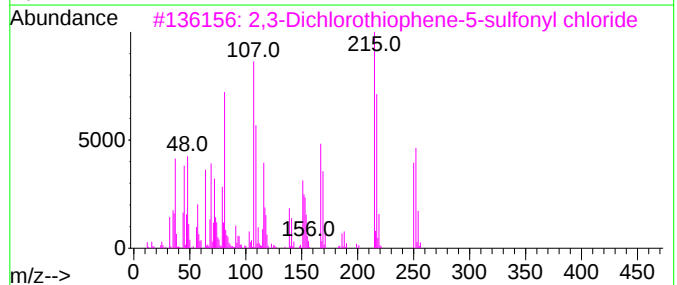
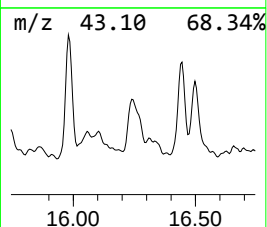
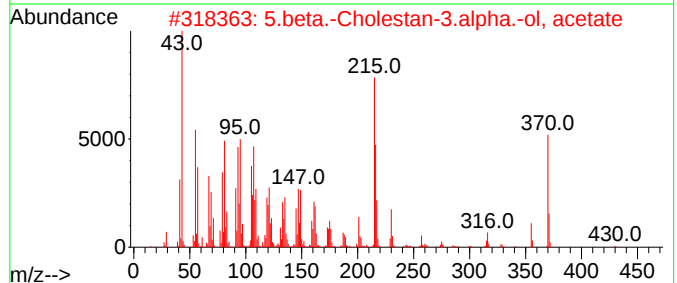
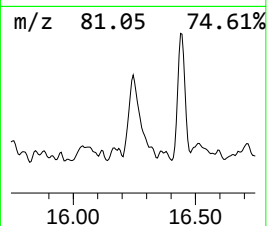
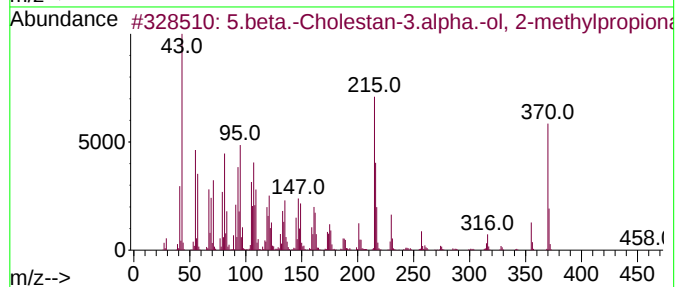
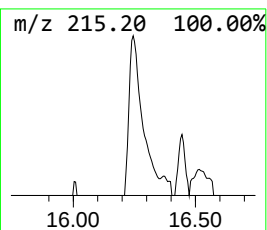
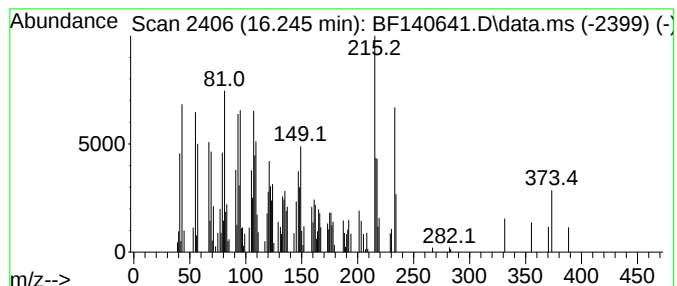
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.245 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.245	6.64 ng	147265	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5.beta.-Cholestan-3.alpha.-ol, 2...	458	C31H54O2	1000514-95-7	49		
2	5.beta.-Cholestan-3.alpha.-ol, a...	430	C29H50O2	033157-47-0	38		
3	2,3-Dichlorothiophene-5-sulfonyl...	250	C4HCl3O2S2	126714-85-0	38		
4	Cholestan-3.beta.-yl benzoate	492	C34H52O2	005808-11-7	27		
5	5-(2-Methyloctan-2-yl)-3-(1R,3R-...	332	C22H36O2	070434-92-3	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
Data File : BF140641.D
Acq On : 26 Nov 2024 14:55
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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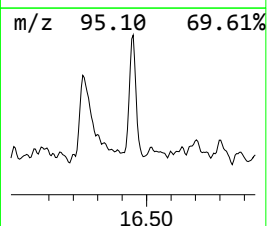
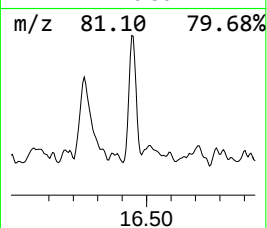
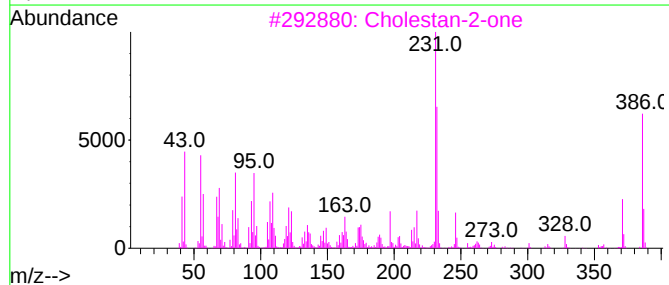
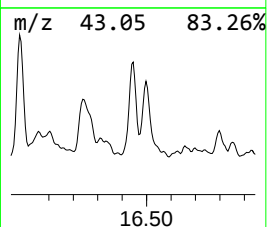
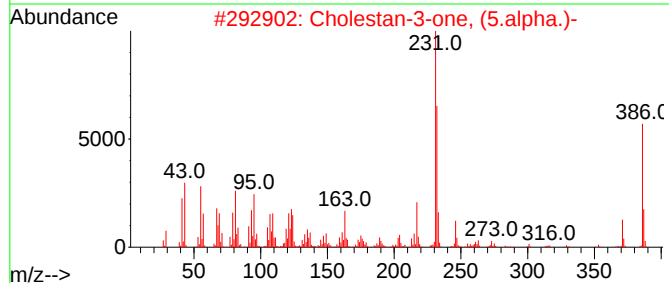
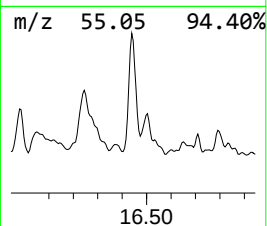
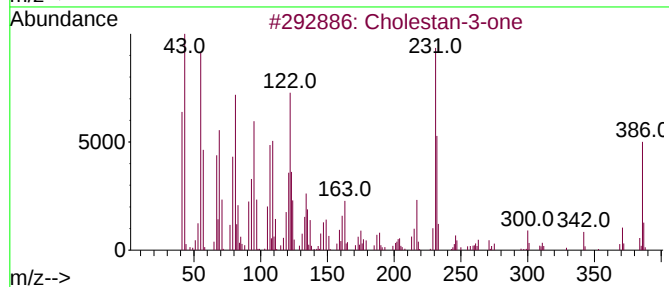
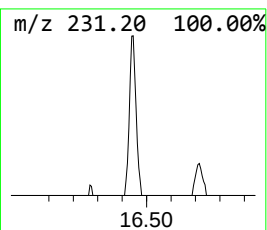
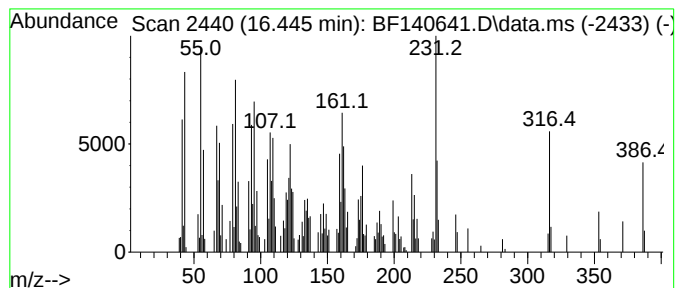
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cholestan-3-one Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.445	6.44 ng	142864	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholestan-3-one	386	C27H46O	015600-08-5	83
2			Cholestan-3-one, (5.alpha.)-	386	C27H46O	000566-88-1	62
3			Cholestan-2-one	386	C27H46O	1010210-43-4	47
4			Cholestan-3-one, (5.beta.)-	386	C27H46O	000601-53-6	42
5			trans-1-(p-Methoxyphenyl)-1-tetr...	316	C21H32O2	123501-67-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
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Acq On : 26 Nov 2024 14:55
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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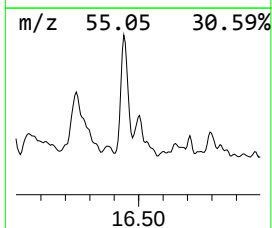
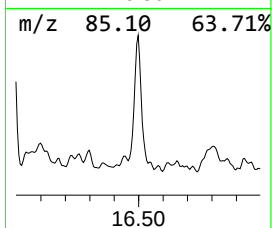
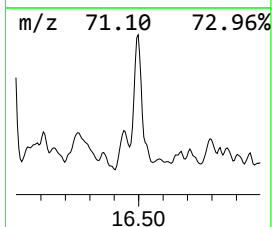
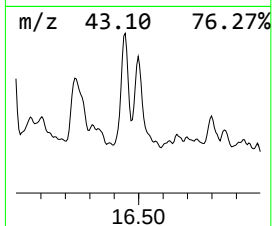
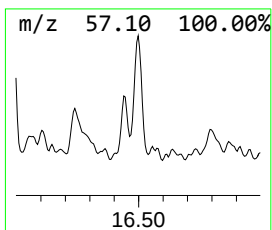
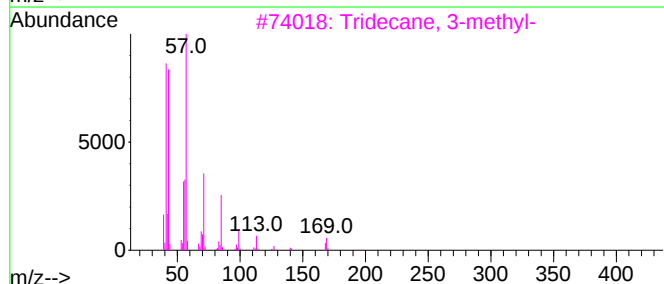
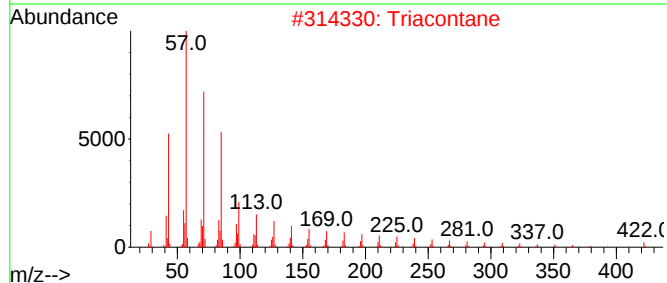
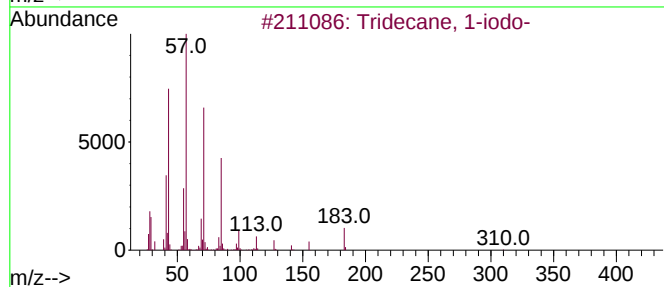
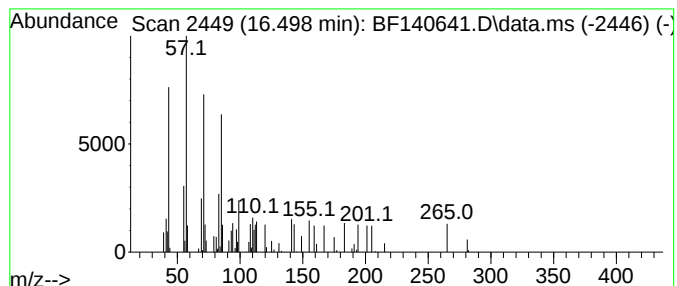
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown16.498 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.498	2.22 ng	49181	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	47
2			Triacotane	422	C30H62	000638-68-6	47
3			Tridecane, 3-methyl-	198	C14H30	006418-41-3	47
4			Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	47
5			Octacosane	394	C28H58	000630-02-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
Data File : BF140641.D
Acq On : 26 Nov 2024 14:55
Operator : RC/JU
Sample : P4860-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PH2-BOT-005

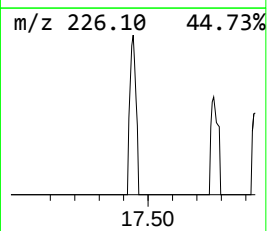
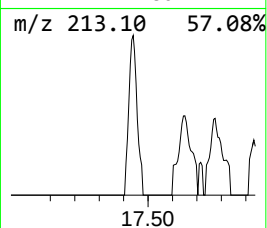
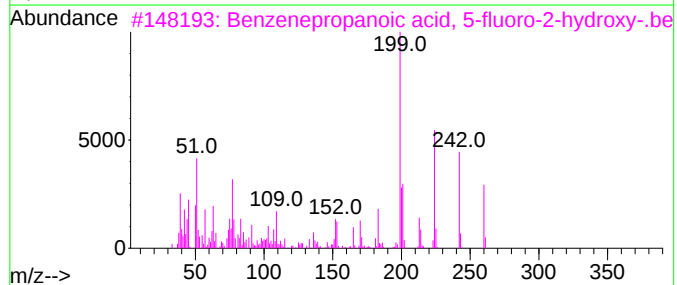
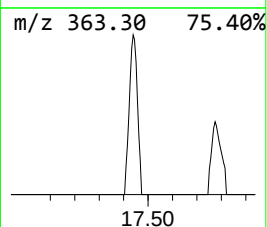
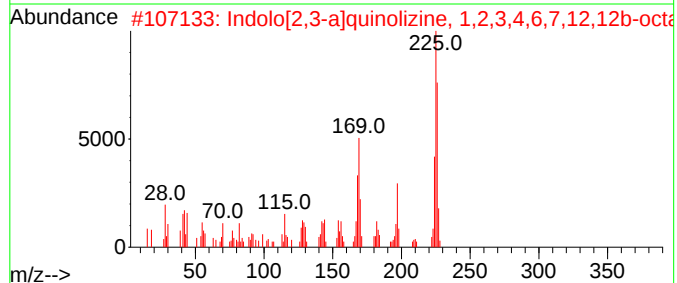
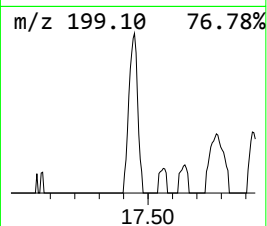
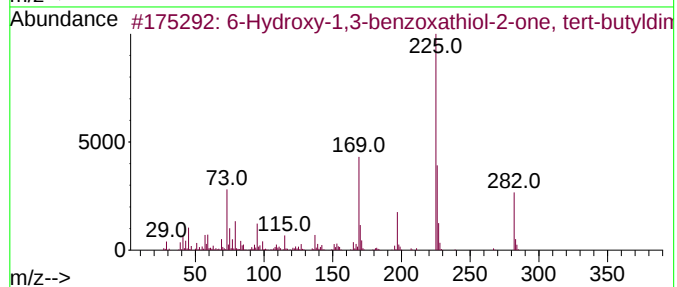
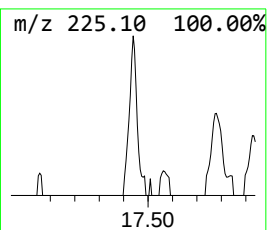
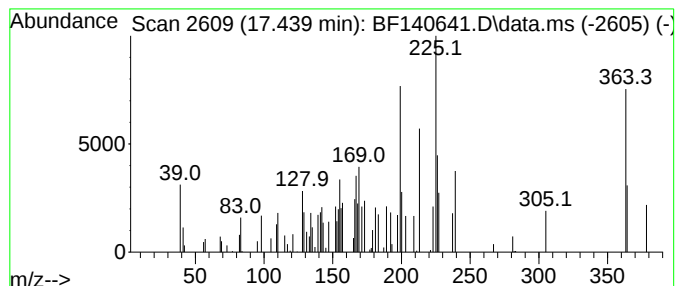
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown17.439 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.439	2.27 ng	50352	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Hydroxy-1,3-benzoxathiol-2-one...	282	C13H18O3SSi	1010461-98-2	12
2			Indolo[2,3-a]quinolizine, 1,2,3,...	226	C15H18N2	004802-79-3	11
3			Benzenepropanoic acid, 5-fluoro-...	260	C15H13FO3	1000362-84-1	10
4			3-Methoxydiphenylamine	199	C13H13NO	000101-16-6	10
5			6-(2-Chloroethylthio)-N,N,N',N'-...	261	C9H16ClN5S	1000102-02-9	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
Data File : BF140641.D
Acq On : 26 Nov 2024 14:55
Operator : RC/JU
Sample : P4860-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PH2-BOT-005

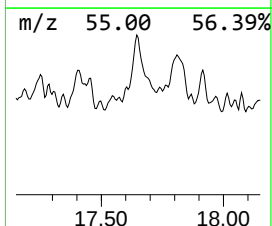
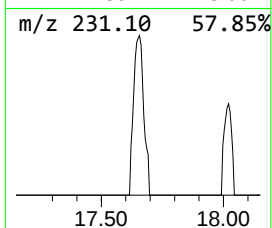
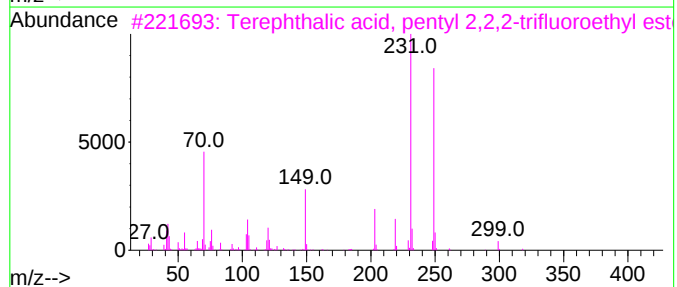
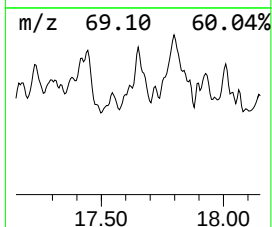
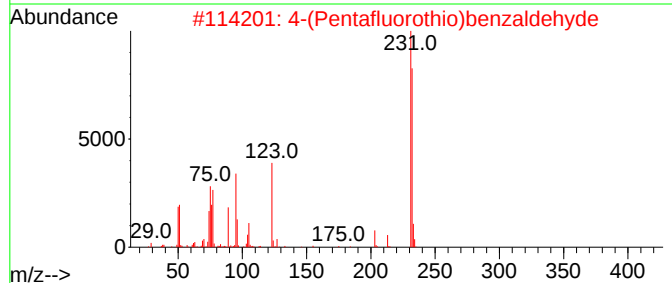
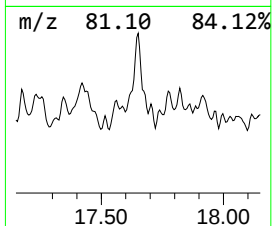
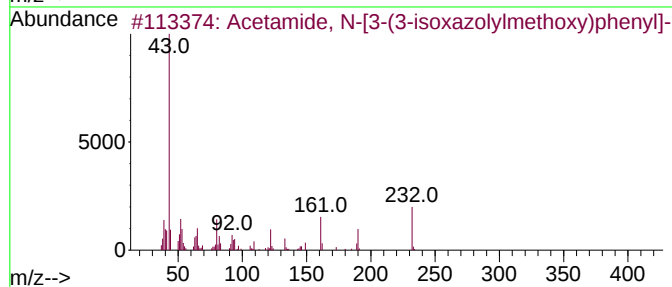
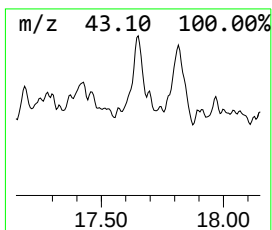
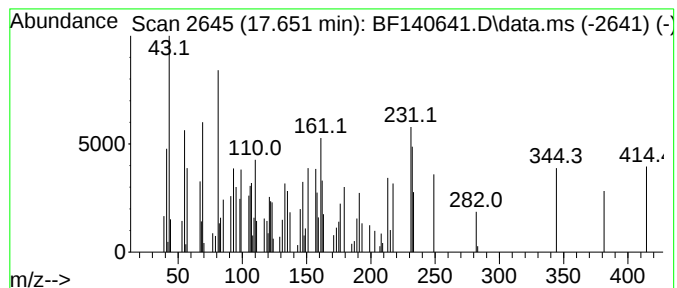
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown17.651 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.651	2.07 ng	45973	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-[3-(3-isoxazolylmet...	232	C12H12N2O3	1000362-73-8	10
2			4-(Pentafluorothio)benzaldehyde	232	C7H5F5OS	401892-84-0	10
3			Terephthalic acid, pentyl 2,2,2-...	318	C15H17F3O4	1000383-04-0	10
4			4-(4-Chlorophenoxy)benzaldehyde	232	C13H9ClO2	061343-99-5	10
5			Costunolide	232	C15H20O2	000553-21-9	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112624\
Data File : BF140641.D
Acq On : 26 Nov 2024 14:55
Operator : RC/JU
Sample : P4860-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PH2-BOT-005

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.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...	2.122	32.7	ng	703888	1	6.869	430326	20.0
Benzophenone	10.616	9.6	ng	284390	3	9.904	590370	20.0
n-Hexadecanoic ...	11.916	4.9	ng	147459	4	11.398	602196	20.0
10,18-Bisnorabi...	12.645	2.4	ng	72175	4	11.398	602196	20.0
Isobutyl tetrac...	13.874	8.3	ng	161630	5	14.045	389191	20.0
Heneicosane	14.192	2.3	ng	45459	5	14.045	389191	20.0
Triacontane	14.504	4.1	ng	80254	5	14.045	389191	20.0
Octacosane, 2-m...	14.815	2.2	ng	48570	6	15.539	443348	20.0
Eicosane	15.157	3.0	ng	65806	6	15.539	443348	20.0
unknown16.245	16.245	6.6	ng	147265	6	15.539	443348	20.0
Cholestan-3-one	16.445	6.4	ng	142864	6	15.539	443348	20.0
unknown16.498	16.498	2.2	ng	49181	6	15.539	443348	20.0
unknown17.439	17.439	2.3	ng	50352	6	15.539	443348	20.0
unknown17.651	17.651	2.1	ng	45973	6	15.539	443348	20.0