

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
Data File : BF140586.D
Acq On : 22 Nov 2024 19:01
Operator : RC/JU
Sample : P4925-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-758

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: BF140586.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	16	rVB	610111	778056	46.34%	6.271%
2	5.499	574	579	594	rVB	1139662	1502757	89.51%	12.111%
3	6.122	680	685	691	rBV5	9445	20312	1.21%	0.164%
4	6.404	729	733	741	rBV4	11899	19982	1.19%	0.161%
5	6.510	745	751	764	rBV	1135088	1556053	92.68%	12.541%
6	6.869	808	812	818	rVB	427545	530561	31.60%	4.276%
7	7.140	852	858	862	rVB4	10049	19824	1.18%	0.160%
8	7.434	903	908	916	rVB	742557	973108	57.96%	7.843%
9	7.504	916	920	924	rVB4	12601	19455	1.16%	0.157%
10	7.681	946	950	956	rVB2	20840	30560	1.82%	0.246%
11	8.151	1025	1030	1036	rBV	503454	641547	38.21%	5.171%
12	8.204	1036	1039	1042	rVV2	22575	30231	1.80%	0.244%
13	8.563	1097	1100	1103	rBV	20854	26908	1.60%	0.217%
14	8.687	1118	1121	1127	rBV4	10556	16813	1.00%	0.136%
15	9.163	1200	1202	1207	rVB2	18823	24214	1.44%	0.195%
16	9.222	1207	1212	1221	rBV	1266374	1678913	100.00%	13.531%
17	9.304	1222	1226	1230	rVB3	15903	22294	1.33%	0.180%
18	9.616	1275	1279	1285	rVB3	20246	29291	1.74%	0.236%
19	9.904	1323	1328	1339	rVB	469976	629925	37.52%	5.077%
20	10.622	1445	1450	1458	rVV	90383	120203	7.16%	0.969%
21	10.698	1458	1463	1476	rVV	615442	818134	48.73%	6.594%
22	10.816	1479	1483	1491	rVB	14508	23350	1.39%	0.188%
23	11.398	1577	1582	1590	rBV	432059	562275	33.49%	4.532%
24	11.916	1666	1670	1684	rBV	75000	145479	8.67%	1.172%
25	12.980	1846	1851	1859	rBV	1025764	1306347	77.81%	10.528%
26	13.874	1999	2003	2015	rBV	44338	93611	5.58%	0.754%
27	14.045	2027	2032	2041	rBV	324051	460121	27.41%	3.708%
28	15.533	2279	2285	2291	rVB	202160	327417	19.50%	2.639%

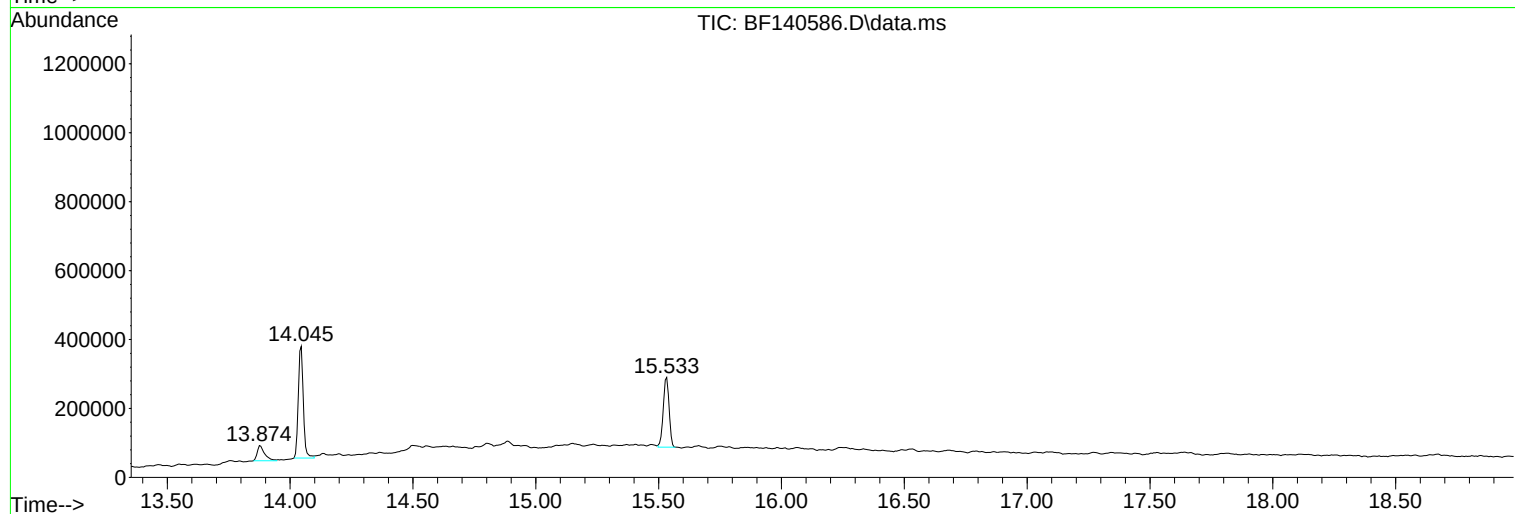
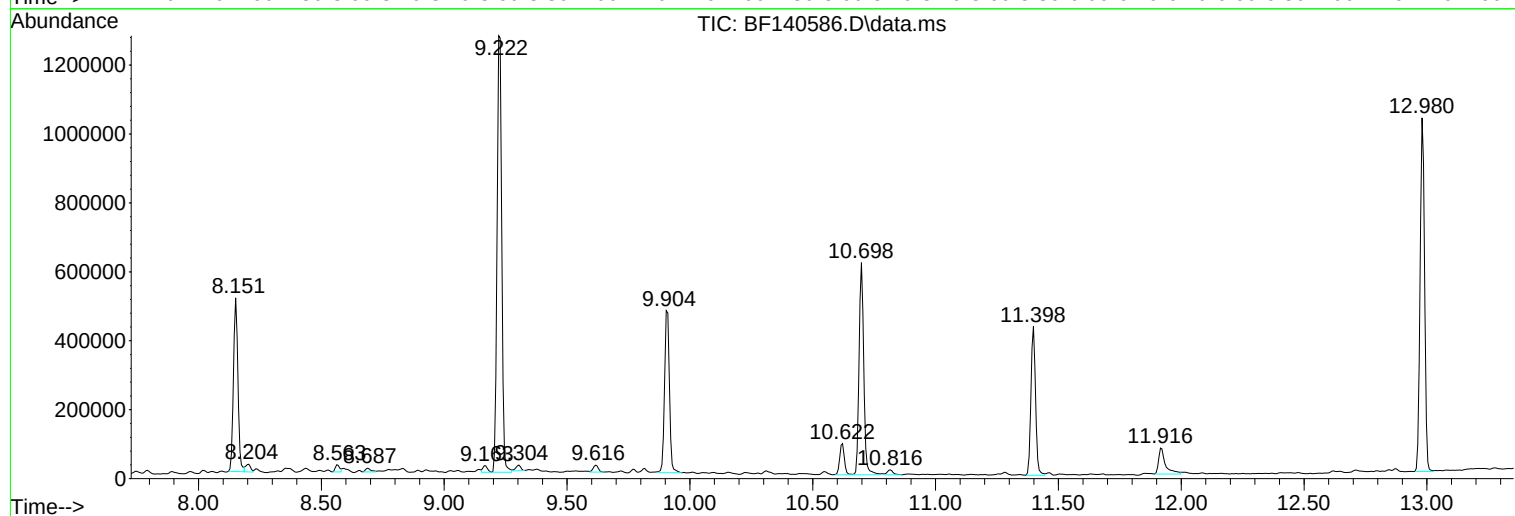
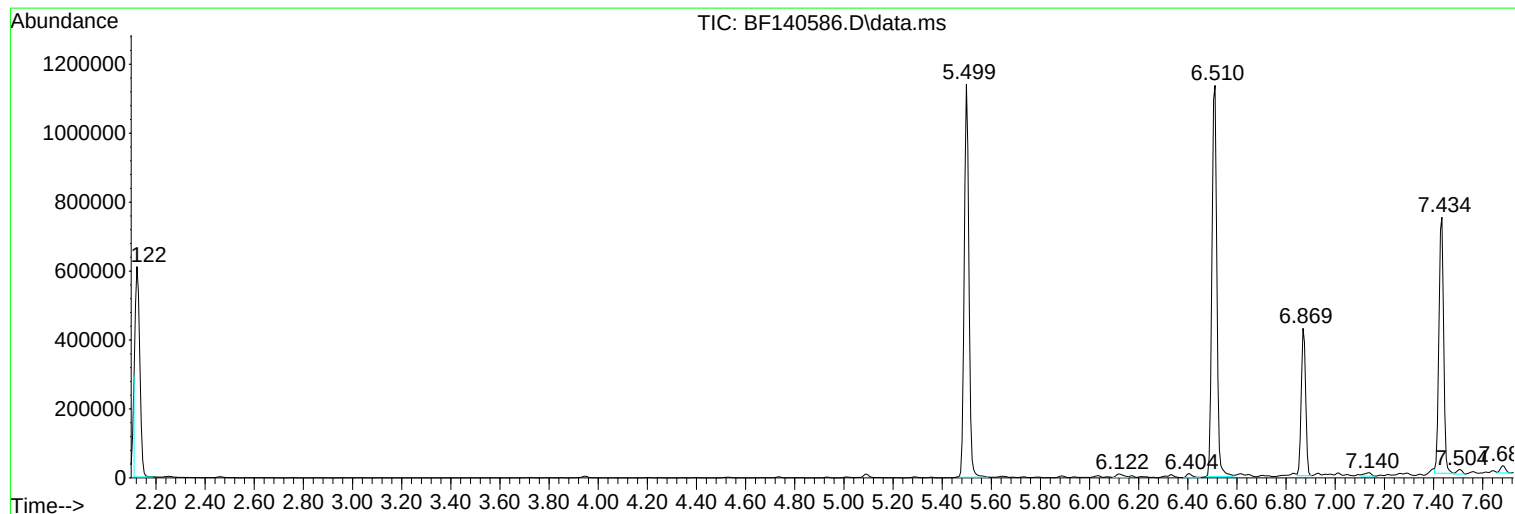
Sum of corrected areas: 12407741

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
Data File : BF140586.D
Acq On : 22 Nov 2024 19:01
Operator : RC/JU
Sample : P4925-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-758

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
Data File : BF140586.D
Acq On : 22 Nov 2024 19:01
Operator : RC/JU
Sample : P4925-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-758

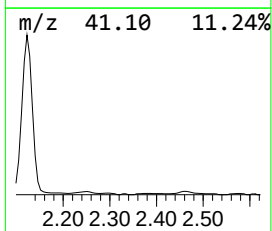
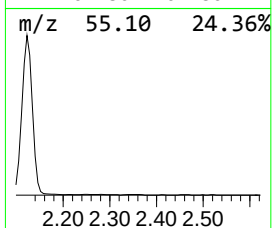
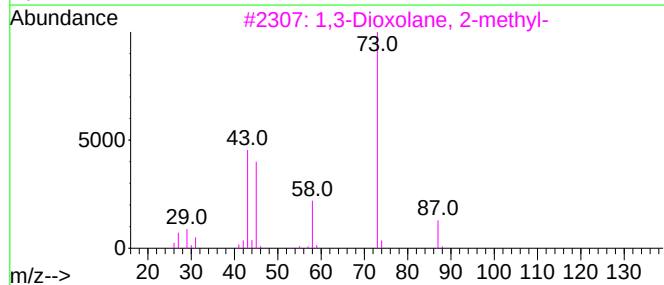
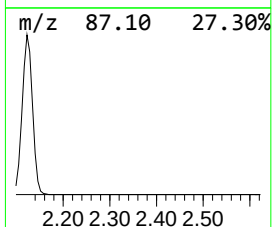
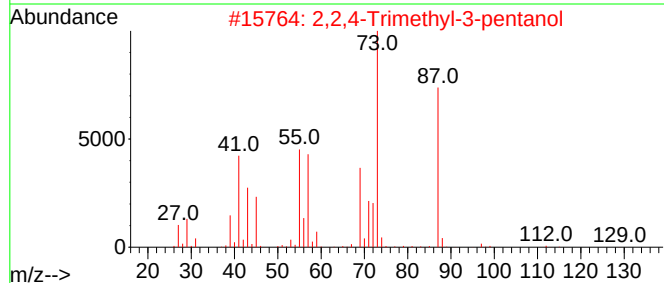
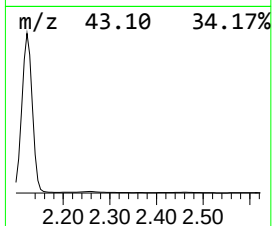
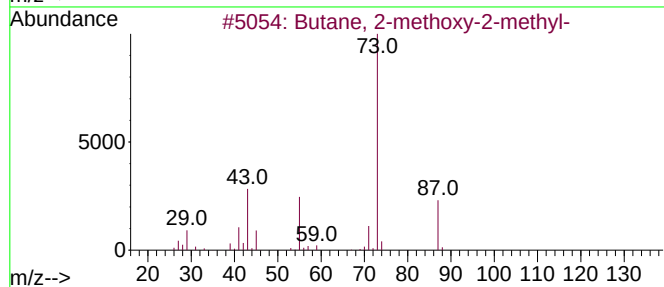
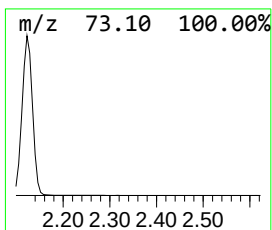
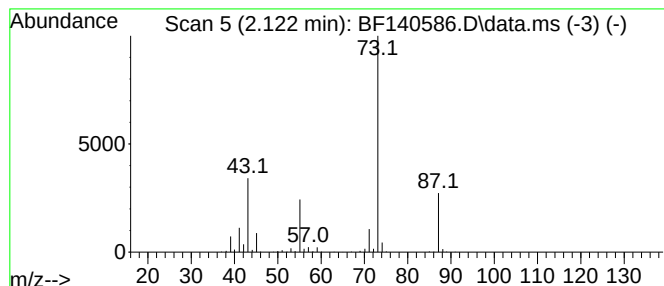
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	29.33 ng	778056	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	78
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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
Data File : BF140586.D
Acq On : 22 Nov 2024 19:01
Operator : RC/JU
Sample : P4925-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-758

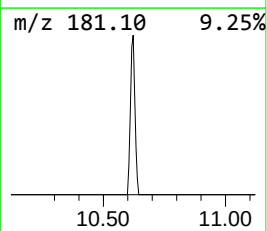
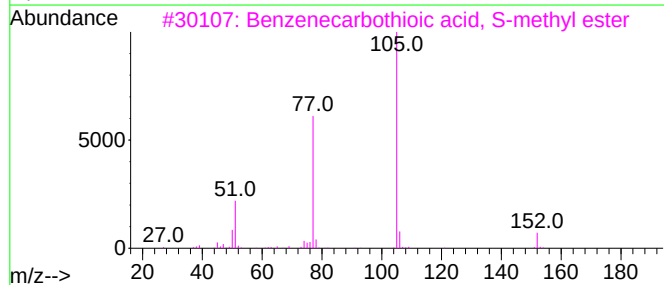
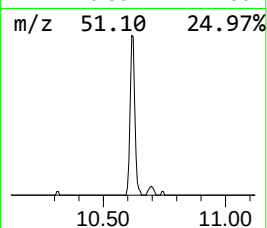
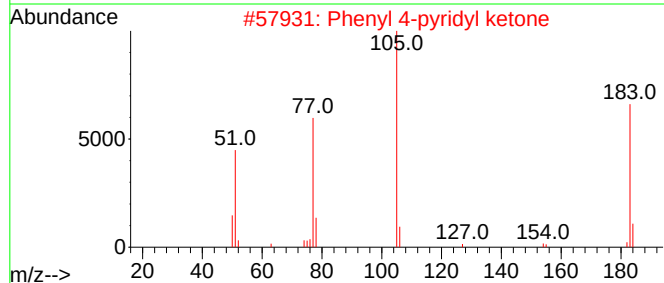
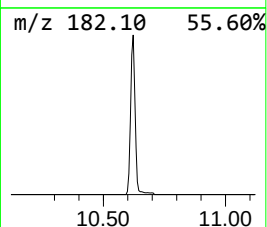
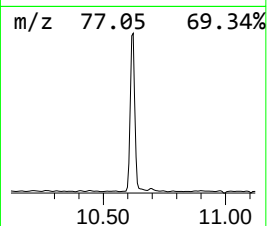
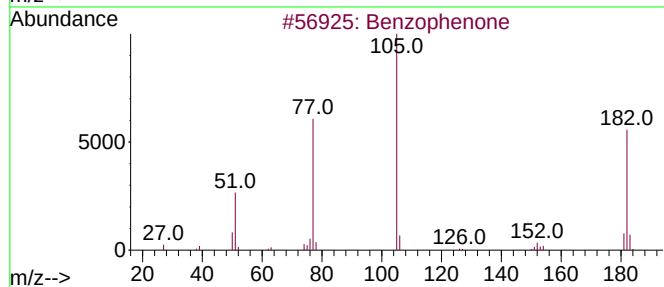
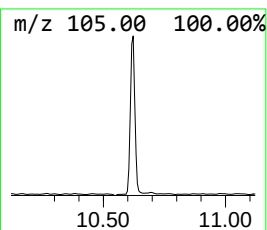
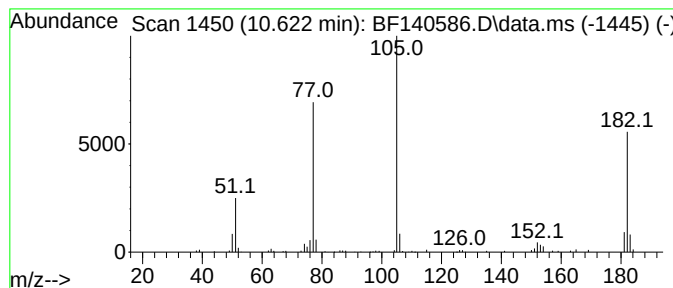
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.622	3.82 ng	120203	Acenaphthene-d10	9.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	53
3			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	53
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	46
5			Benzenecarbothioic acid	138	C7H6OS	000098-91-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
Data File : BF140586.D
Acq On : 22 Nov 2024 19:01
Operator : RC/JU
Sample : P4925-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-758

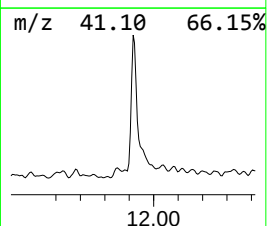
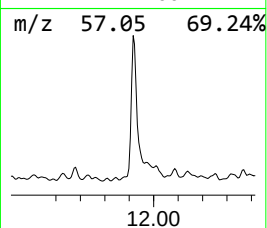
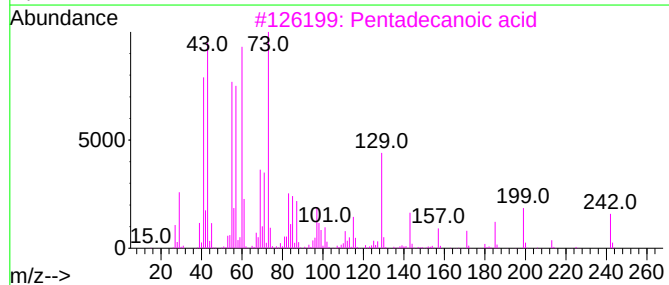
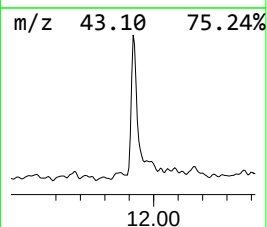
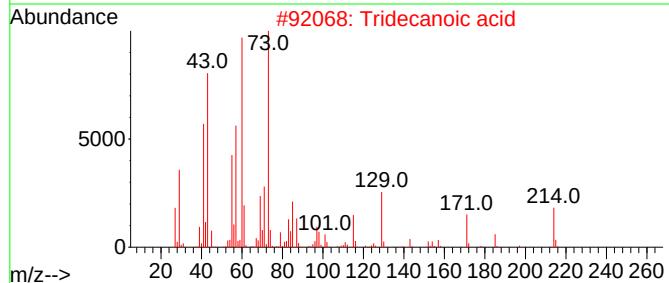
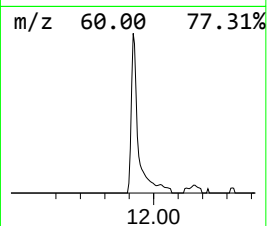
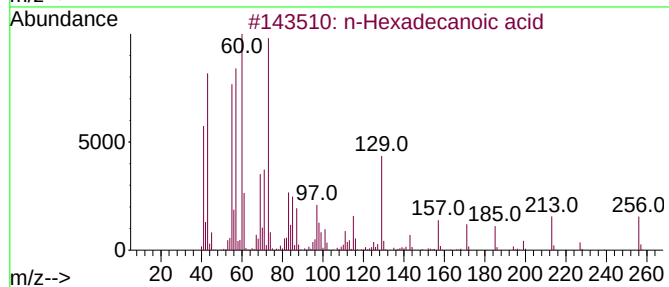
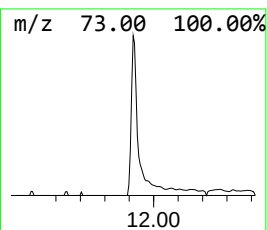
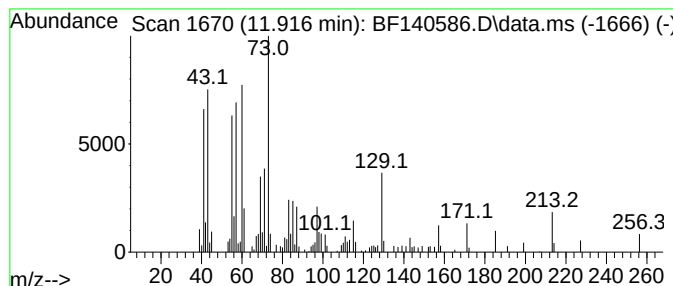
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	5.17 ng	145479	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	93
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5			n-Decanoic acid	172	C10H20O2	000334-48-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
Data File : BF140586.D
Acq On : 22 Nov 2024 19:01
Operator : RC/JU
Sample : P4925-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-758

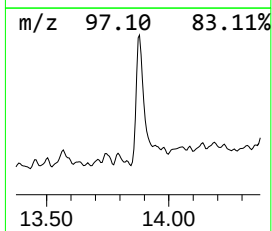
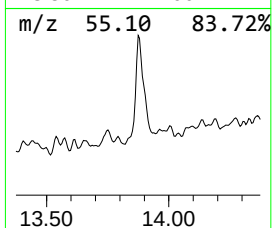
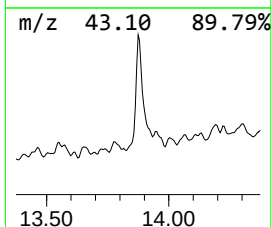
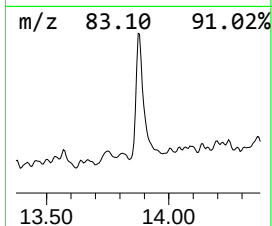
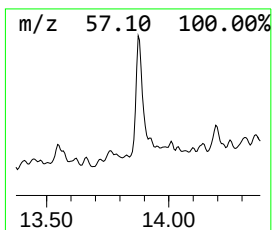
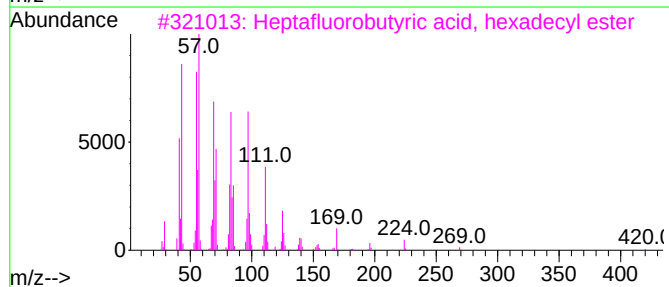
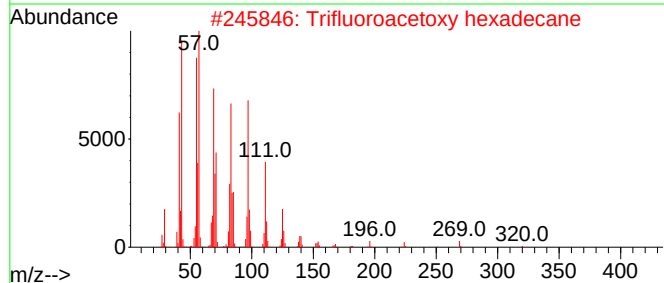
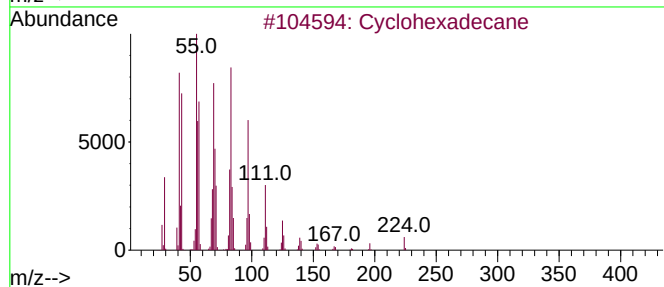
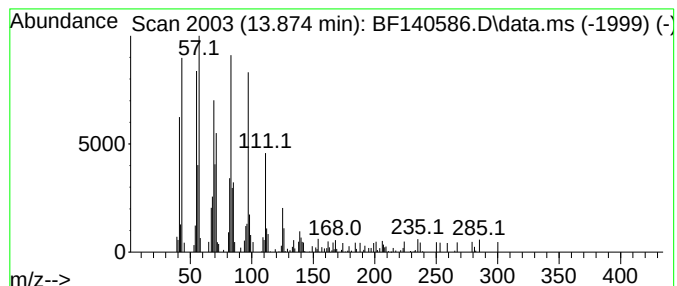
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Cyclohexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	4.07 ng	93611	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	95
2			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93
3			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5			Octacosanol	410	C28H58O	000557-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
Data File : BF140586.D
Acq On : 22 Nov 2024 19:01
Operator : RC/JU
Sample : P4925-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-758

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.122	29.3	ng	778056	1	6.869	530561	20.0
Benzophenone	10.622	3.8	ng	120203	3	9.910	629925	20.0
n-Hexadecanoic ...	11.916	5.2	ng	145479	4	11.398	562275	20.0
Cyclohexadecane	13.874	4.1	ng	93611	5	14.045	460121	20.0