

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
 Data File : BF139988.D
 Acq On : 24 Oct 2024 02:34
 Operator : RC/JU
 Sample : P4468-03
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-329

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-BF101824.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139988.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.193	8	17	24	rBV	1680825	2708767	100.00%	12.575%
2	5.116	504	514	519	rBV	33674	53407	1.97%	0.248%
3	5.504	574	580	586	rBV	1727110	2315237	85.47%	10.748%
4	6.510	744	751	763	rBV	1636848	2236679	82.57%	10.383%
5	6.892	810	816	821	rVB	574248	713845	26.35%	3.314%
6	7.445	904	910	915	rBV	1061233	1393345	51.44%	6.468%
7	7.698	949	953	959	rBV	22843	28832	1.06%	0.134%
8	8.169	1028	1033	1042	rBV	657625	838019	30.94%	3.890%
9	9.245	1210	1216	1226	rBV	1762546	2216628	81.83%	10.290%
10	9.927	1326	1332	1342	rVB	650793	832772	30.74%	3.866%
11	10.639	1447	1453	1460	rBV	388047	484016	17.87%	2.247%
12	10.710	1460	1465	1481	rBV	1023049	1342299	49.55%	6.231%
13	10.945	1495	1505	1509	rBV	45800	63407	2.34%	0.294%
14	11.410	1579	1584	1593	rBV2	649385	1017049	37.55%	4.721%
15	11.486	1593	1597	1601	rVB	36343	48999	1.81%	0.227%
16	11.639	1618	1623	1630	rBV2	20201	34633	1.28%	0.161%
17	11.933	1665	1673	1685	rBV2	289114	506651	18.70%	2.352%
18	12.027	1685	1689	1697	rVV3	39244	76838	2.84%	0.357%
19	12.204	1715	1719	1727	rBV3	14476	34557	1.28%	0.160%
20	12.439	1755	1759	1766	rVB2	18503	30519	1.13%	0.142%
21	12.627	1786	1791	1795	rBV	155876	213723	7.89%	0.992%
22	12.721	1803	1807	1812	rBV	67760	105394	3.89%	0.489%
23	12.851	1820	1829	1834	rBV	124415	222937	8.23%	1.035%
24	12.998	1848	1854	1862	rBV	1884774	2385363	88.06%	11.073%
25	13.898	2003	2007	2018	rBV2	118007	217493	8.03%	1.010%
26	14.051	2028	2033	2045	rVB	615912	911163	33.64%	4.230%
27	15.533	2280	2285	2297	rVB	310874	509164	18.80%	2.364%

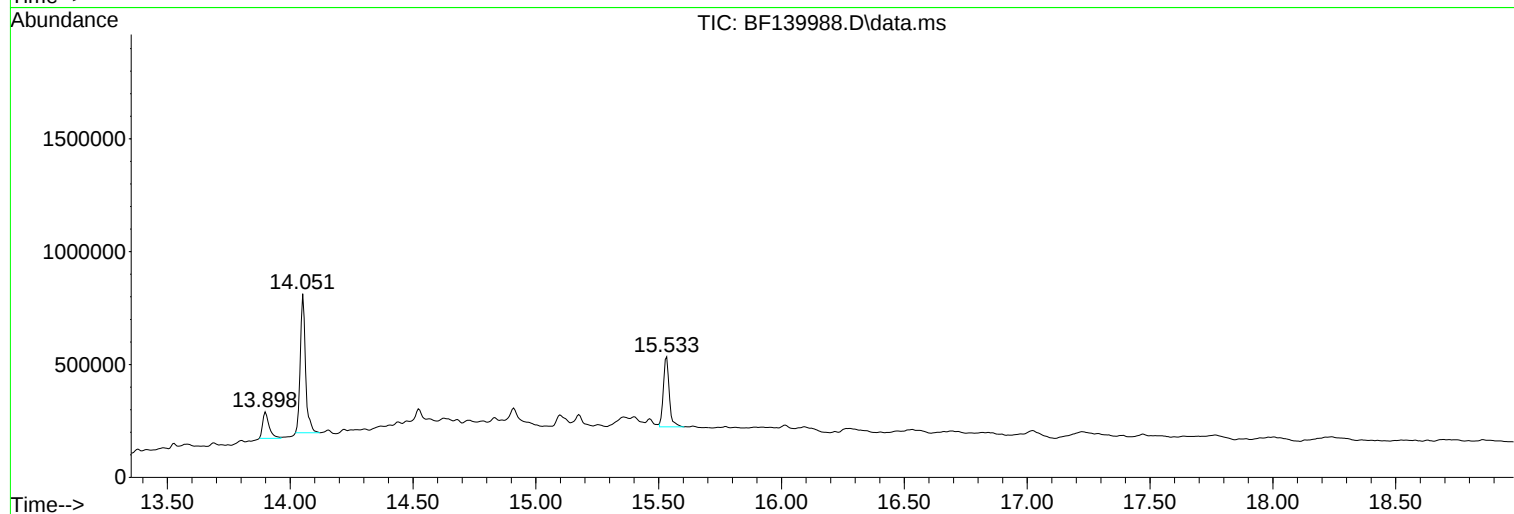
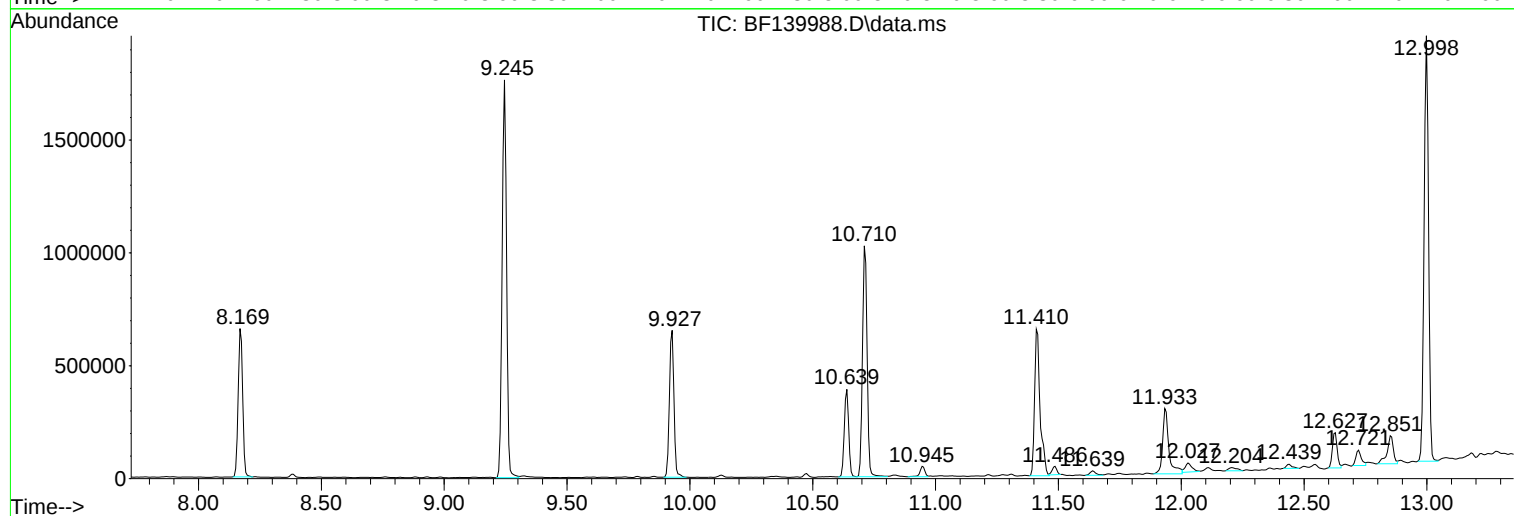
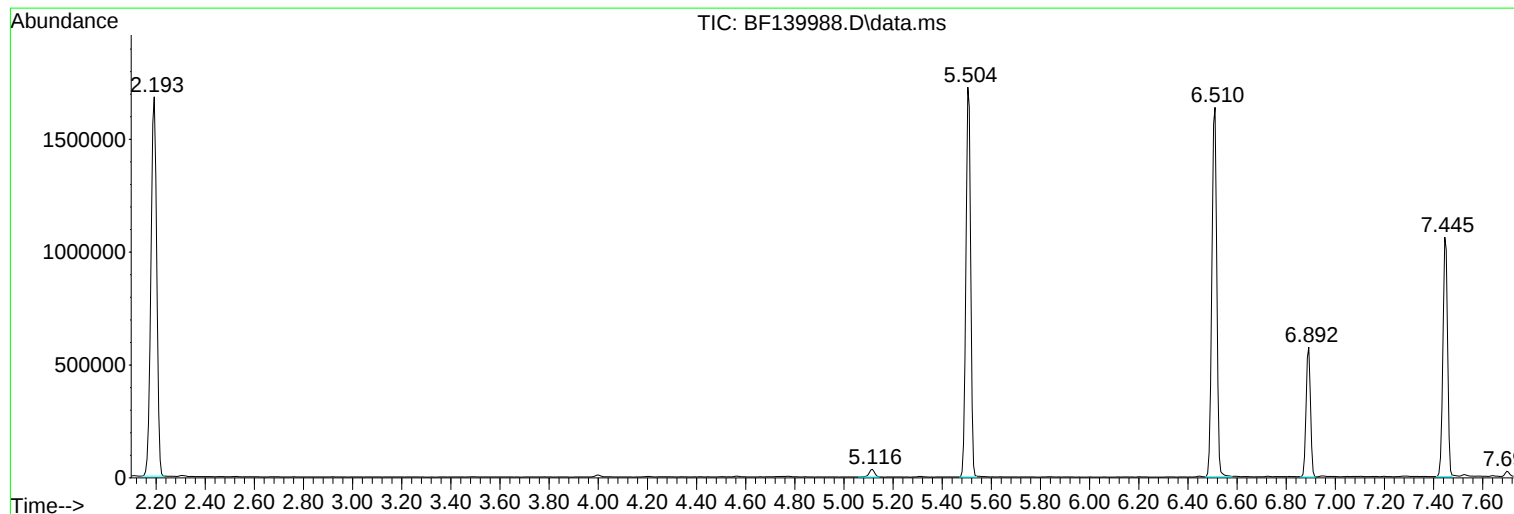
Sum of corrected areas: 21541736

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
Data File : BF139988.D
Acq On : 24 Oct 2024 02:34
Operator : RC/JU
Sample : P4468-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-329

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.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\BF102324\
Data File : BF139988.D
Acq On : 24 Oct 2024 02:34
Operator : RC/JU
Sample : P4468-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-329

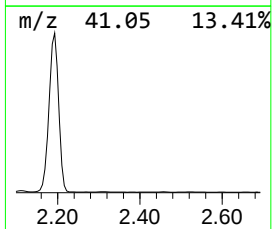
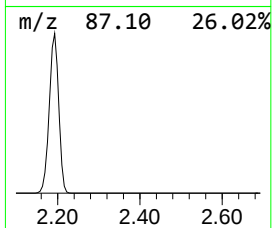
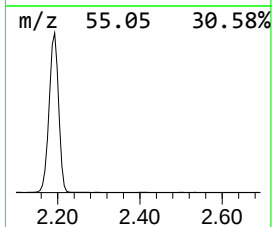
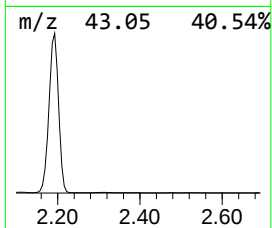
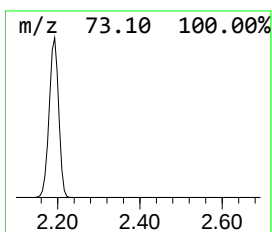
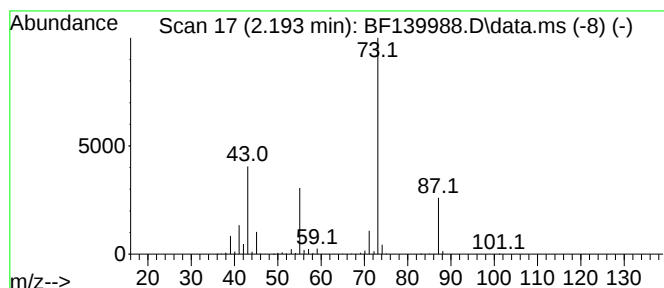
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.193	75.89 ng	2708770	1,4-Dichlorobenzene-d4	6.892

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	12
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
Data File : BF139988.D
Acq On : 24 Oct 2024 02:34
Operator : RC/JU
Sample : P4468-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-329

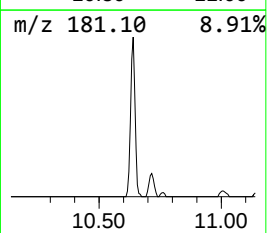
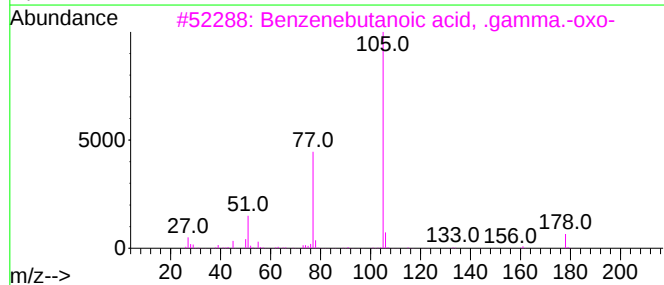
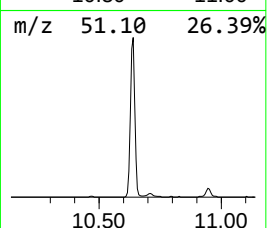
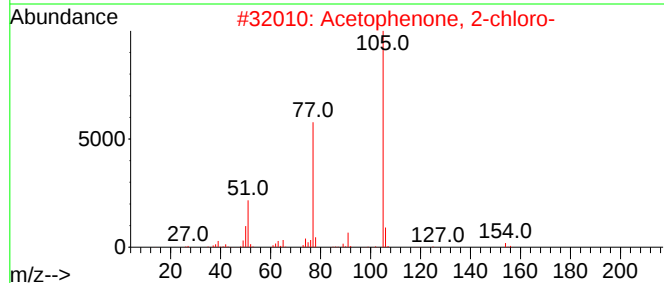
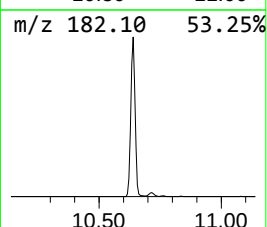
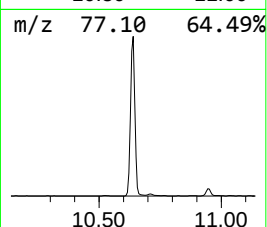
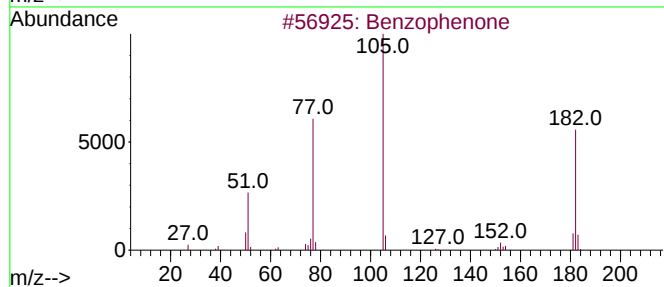
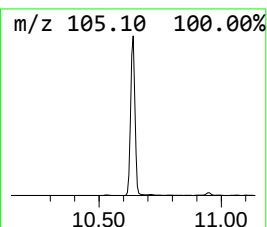
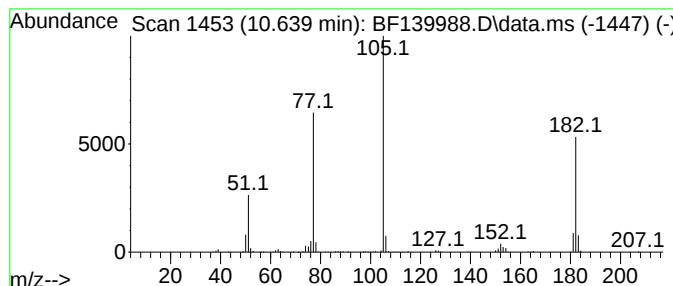
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.639	11.62 ng	484016	Acenaphthene-d10	9.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
3			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47
4			Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	47
5			Vinyl benzoate	148	C9H8O2	000769-78-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
Data File : BF139988.D
Acq On : 24 Oct 2024 02:34
Operator : RC/JU
Sample : P4468-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-329

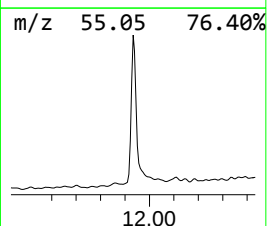
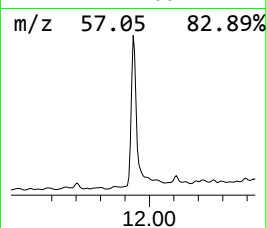
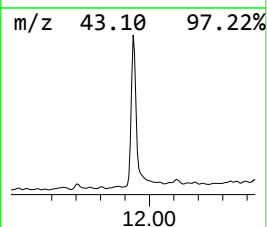
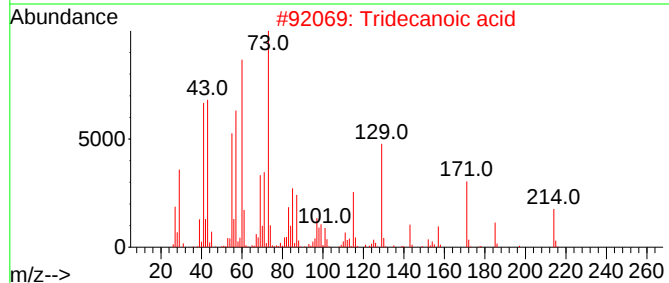
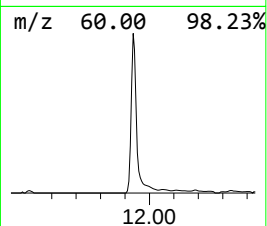
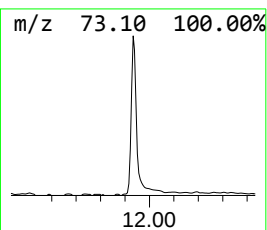
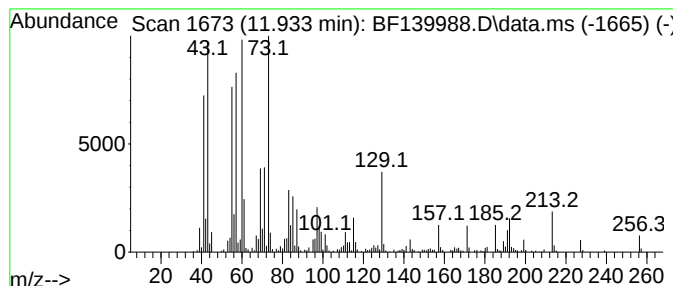
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.933	9.96 ng	506651	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	95
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			n-Decanoic acid	172	C10H20O2	000334-48-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
Data File : BF139988.D
Acq On : 24 Oct 2024 02:34
Operator : RC/JU
Sample : P4468-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-329

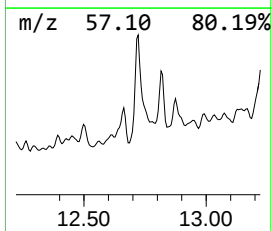
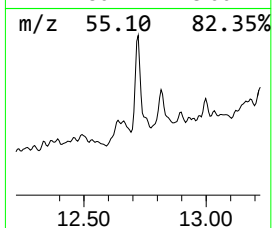
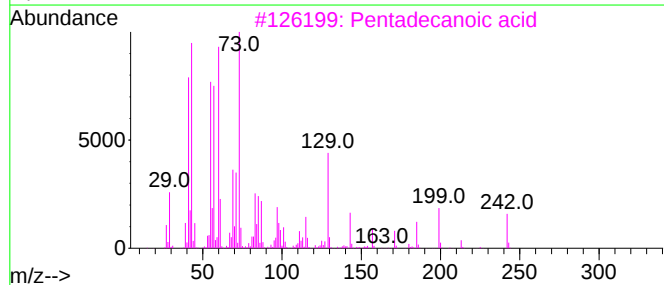
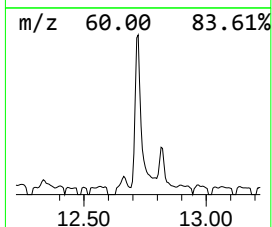
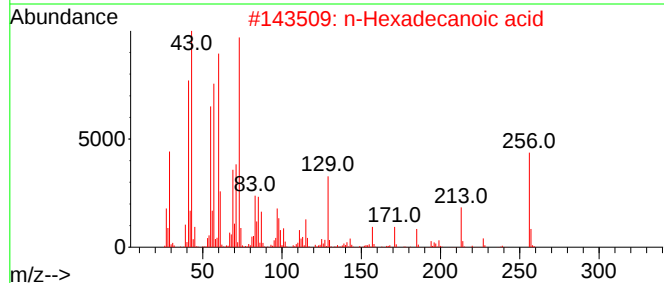
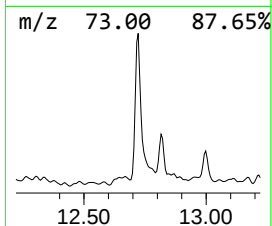
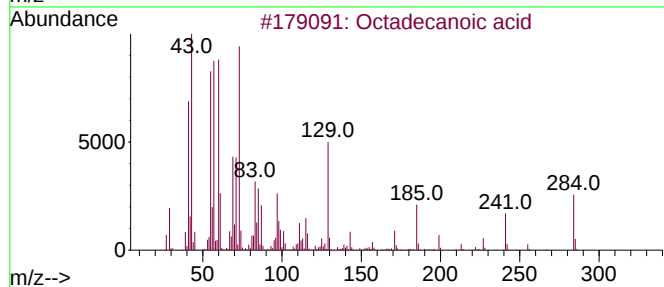
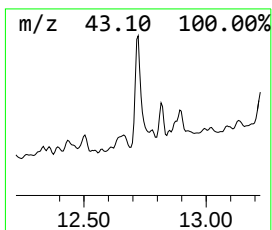
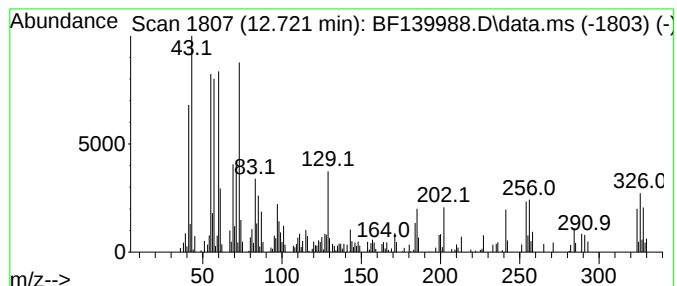
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.721	2.07 ng	105394	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	60
4			Heneicosanoic acid	326	C21H42O2	002363-71-5	56
5			Dodecanoic acid	200	C12H24O2	000143-07-7	48



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
Data File : BF139988.D
Acq On : 24 Oct 2024 02:34
Operator : RC/JU
Sample : P4468-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-329

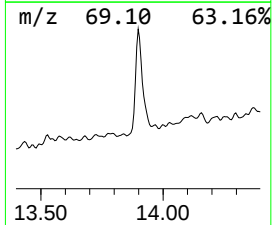
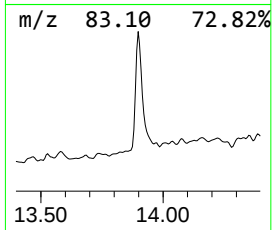
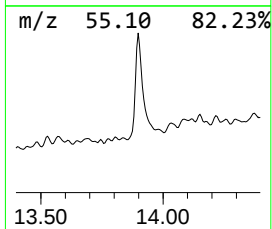
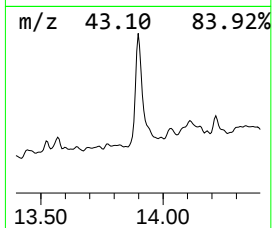
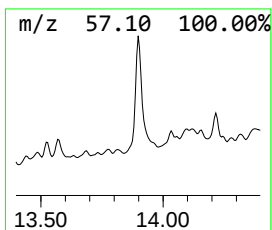
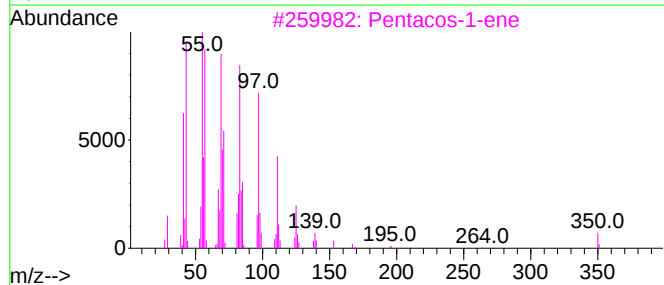
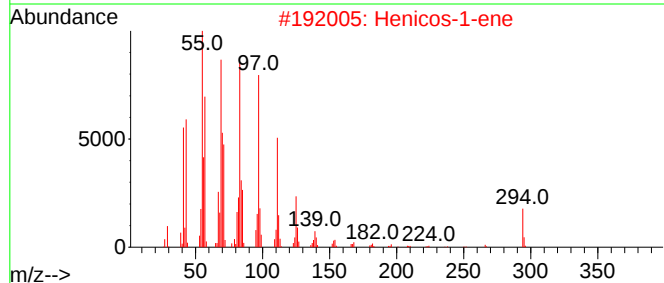
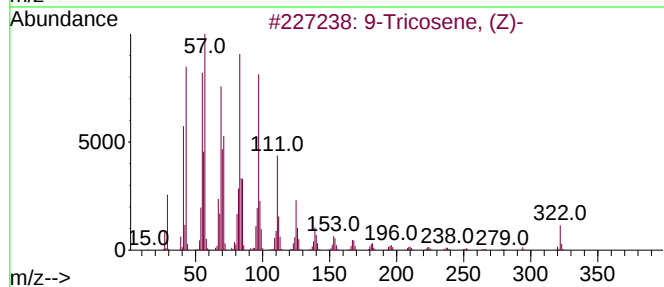
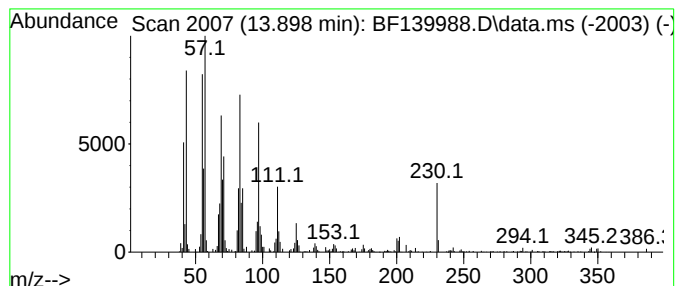
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 9-Tricosene, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.898	4.77 ng	217493	Chrysene-d12	14.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Tricosene, (Z)-	322	C23H46	027519-02-4	99
2			Henicos-1-ene	294	C21H42	001599-68-4	98
3			Pentacos-1-ene	350	C25H50	016980-85-1	94
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
5			1-Hexacosene	364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
Data File : BF139988.D
Acq On : 24 Oct 2024 02:34
Operator : RC/JU
Sample : P4468-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-329

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.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...	2.193	75.9	ng	2708770	1	6.892	713845	20.0
Benzophenone	10.639	11.6	ng	484016	3	9.927	832772	20.0
n-Hexadecanoic ...	11.933	10.0	ng	506651	4	11.410	1017050	20.0
Octadecanoic acid	12.721	2.1	ng	105394	4	11.410	1017050	20.0
9-Tricosene, (Z)-	13.898	4.8	ng	217493	5	14.051	911163	20.0