

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
 Data File : BF140606.D
 Acq On : 25 Nov 2024 16:49
 Operator : RC/JU
 Sample : P4947-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A3988

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: BF140606.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	19	rVB	1976303	2364343	76.37%	12.108%
2	5.087	504	509	516	rVB	23719	36354	1.17%	0.186%
3	5.498	573	579	594	rVB	908550	1250006	40.37%	6.401%
4	6.504	743	750	765	rBV	605602	828723	26.77%	4.244%
5	6.869	807	812	819	rBV	342866	424929	13.73%	2.176%
6	7.434	899	908	913	rBV	1145492	1557642	50.31%	7.977%
7	7.681	944	950	957	rVB2	20691	35690	1.15%	0.183%
8	7.887	978	985	989	rBV2	31764	50552	1.63%	0.259%
9	8.069	1011	1016	1023	rBV6	13664	33063	1.07%	0.169%
10	8.151	1024	1030	1041	rVB	420577	585278	18.90%	2.997%
11	8.475	1081	1085	1092	rBV3	20999	51327	1.66%	0.263%
12	8.651	1111	1115	1120	rVB2	39736	51274	1.66%	0.263%
13	9.222	1207	1212	1225	rVB	2281291	2996429	96.78%	15.345%
14	9.563	1265	1270	1278	rVB6	16266	41397	1.34%	0.212%
15	9.904	1323	1328	1337	rVB	496401	625933	20.22%	3.205%
16	10.316	1393	1398	1405	rVB2	20316	38196	1.23%	0.196%
17	10.616	1444	1449	1454	rBV	205972	276390	8.93%	1.415%
18	10.698	1457	1463	1478	rVV	1457521	1980049	63.95%	10.140%
19	10.810	1479	1482	1490	rVB5	14668	31267	1.01%	0.160%
20	10.928	1498	1502	1510	rVB	44908	61363	1.98%	0.314%
21	11.251	1553	1557	1564	rVB2	51132	77569	2.51%	0.397%
22	11.392	1576	1581	1588	rBV	495989	642941	20.77%	3.292%
23	11.775	1641	1646	1653	rBV	119835	166454	5.38%	0.852%
24	11.916	1664	1670	1680	rBV2	263300	417333	13.48%	2.137%
25	12.039	1687	1691	1695	rVV	22800	33961	1.10%	0.174%
26	12.610	1779	1788	1799	rBV2	48942	148055	4.78%	0.758%
27	12.704	1799	1804	1814	rBV2	92483	156964	5.07%	0.804%
28	12.980	1845	1851	1857	rBV	2295511	3096011	100.00%	15.855%
29	13.410	1918	1924	1934	rVB2	43298	100242	3.24%	0.513%
30	13.904	2002	2008	2023	rBV2	45857	115201	3.72%	0.590%
31	14.045	2024	2032	2040	rBV	391820	567146	18.32%	2.904%
32	14.574	2116	2122	2132	rVB	45136	83126	2.68%	0.426%
33	14.898	2172	2177	2182	rVB	44519	63491	2.05%	0.325%
34	15.080	2203	2208	2214	rBV	33523	59931	1.94%	0.307%
35	15.539	2280	2286	2296	rBV	244486	392435	12.68%	2.010%
36	15.651	2300	2305	2313	rVB	22048	45179	1.46%	0.231%

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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

37	17.474	2609	2615	2628	rVB4	15384	41374	1.34%	0.212%
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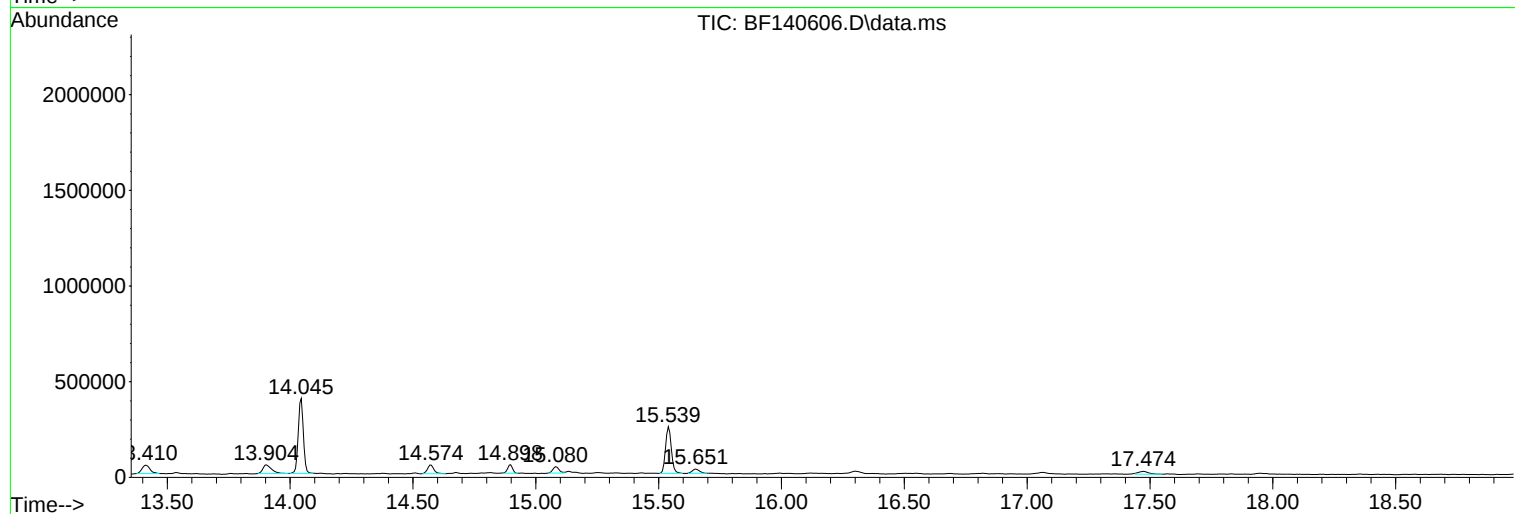
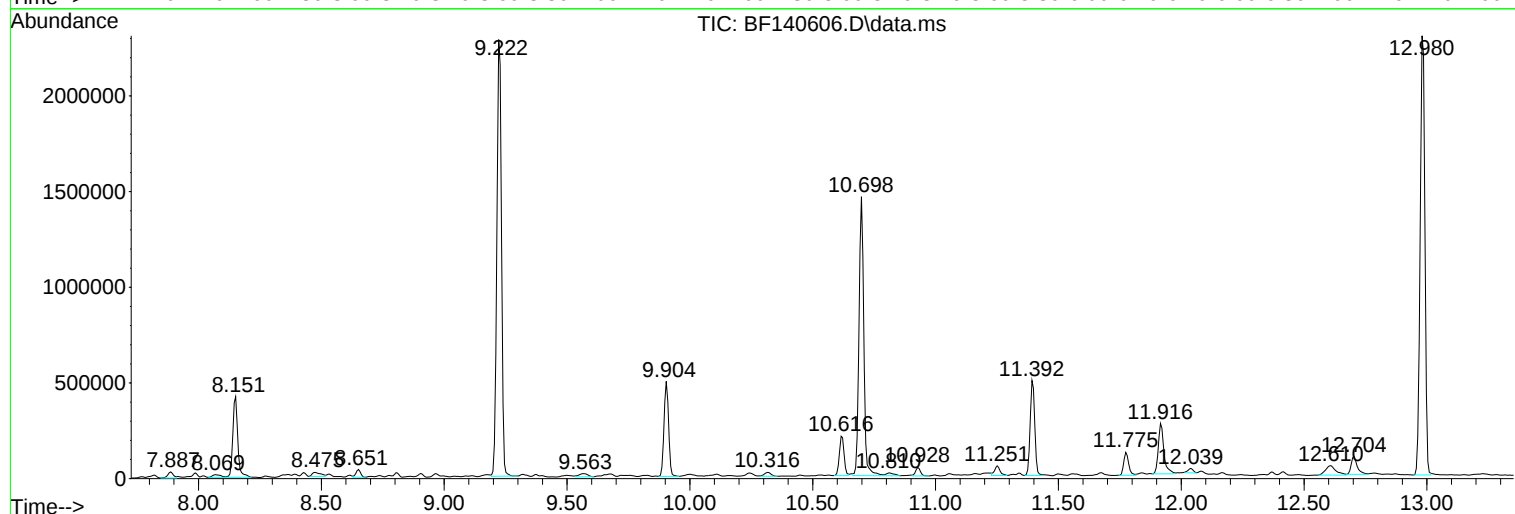
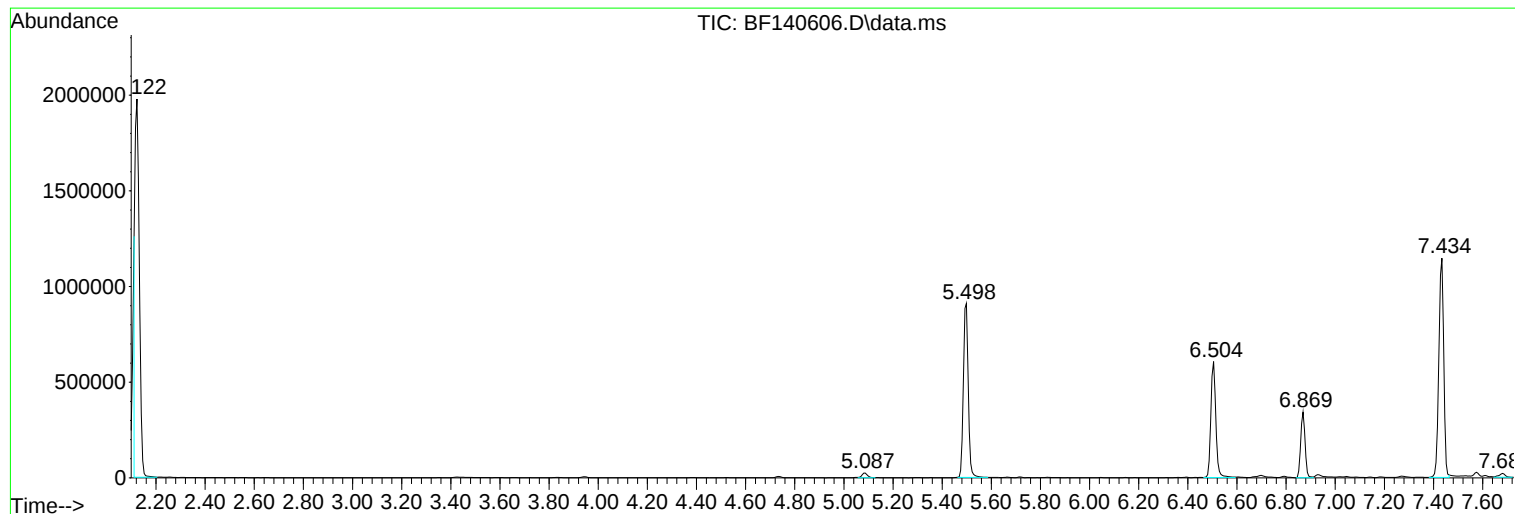
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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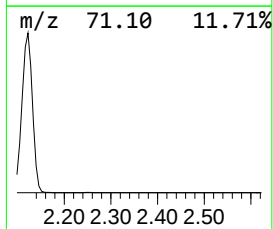
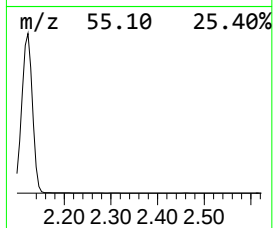
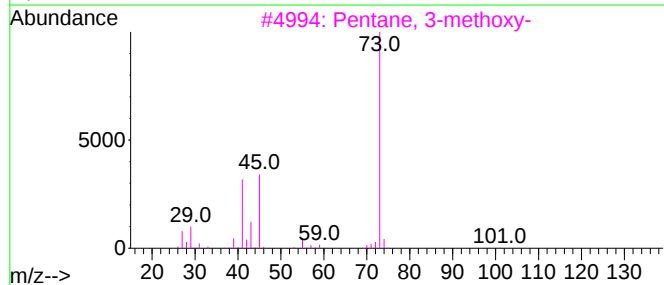
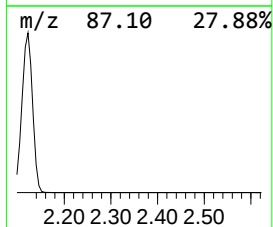
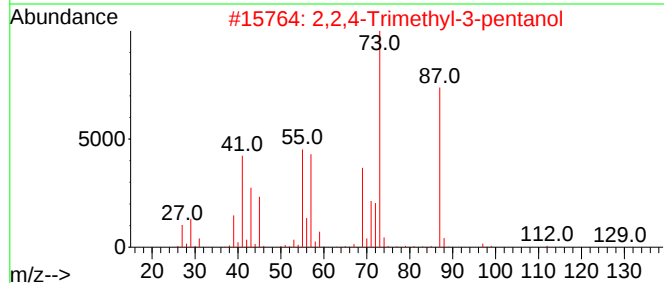
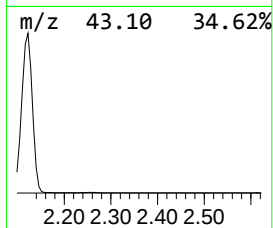
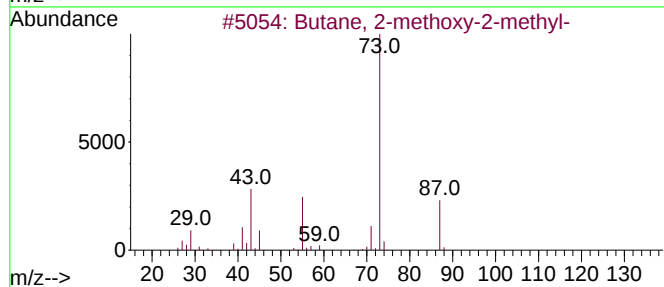
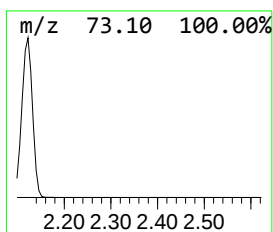
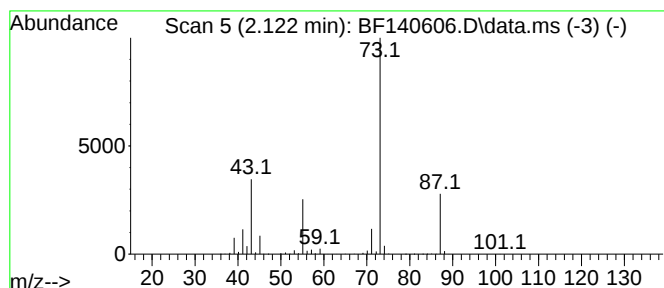
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	111.28 ng	2364340	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			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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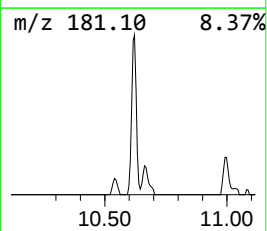
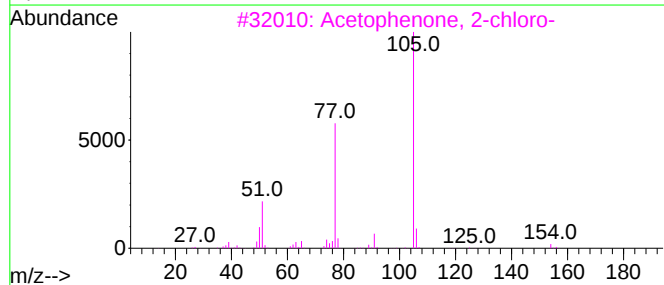
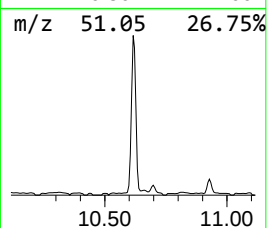
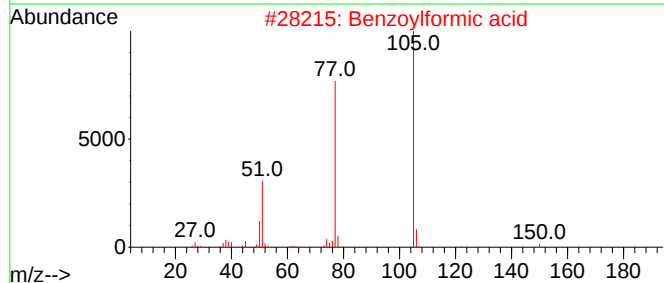
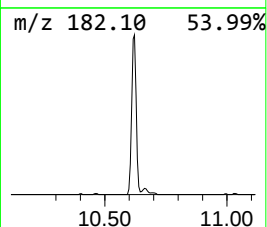
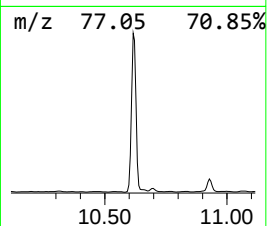
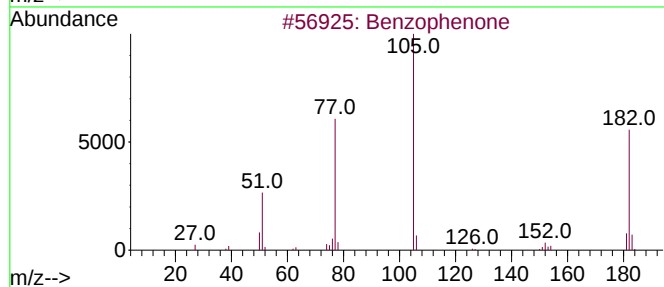
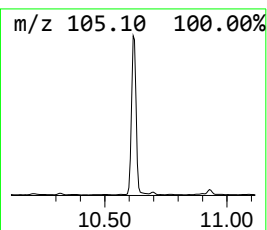
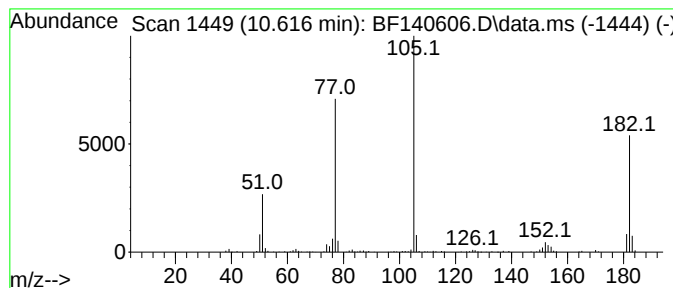
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.616	8.83 ng	276390	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzoylformic acid	150	C8H6O3	000611-73-4	53
3			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	49
5			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	47



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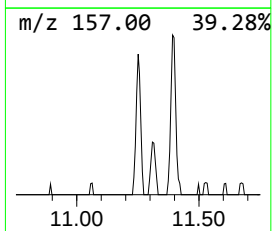
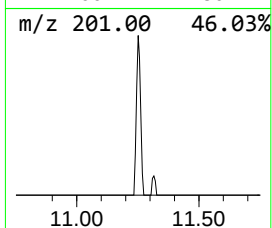
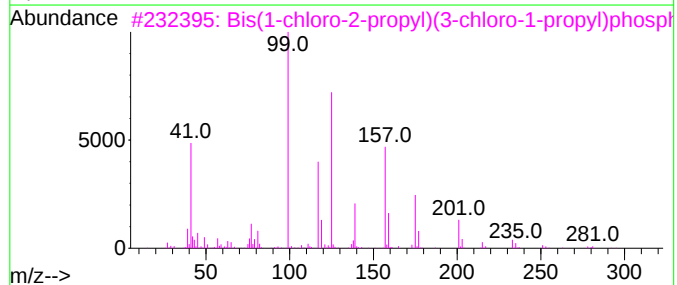
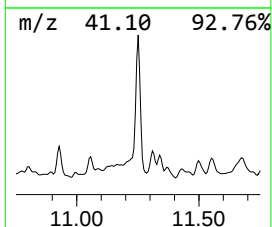
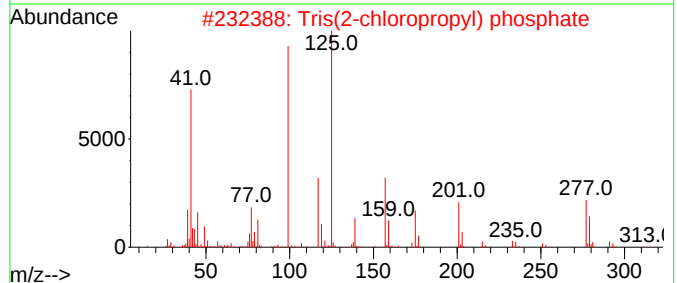
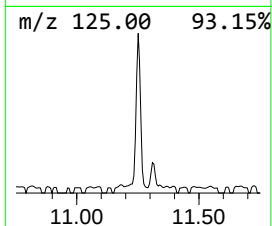
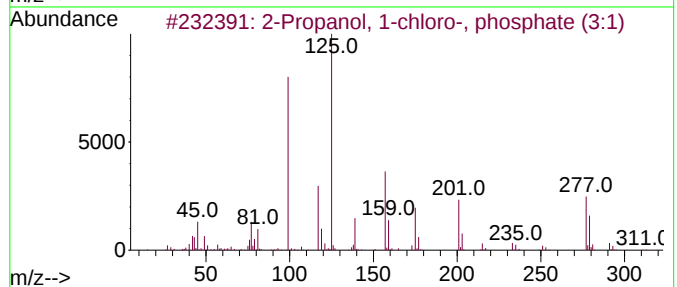
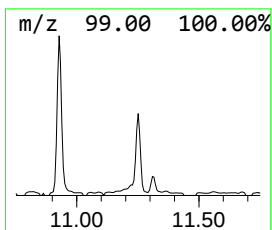
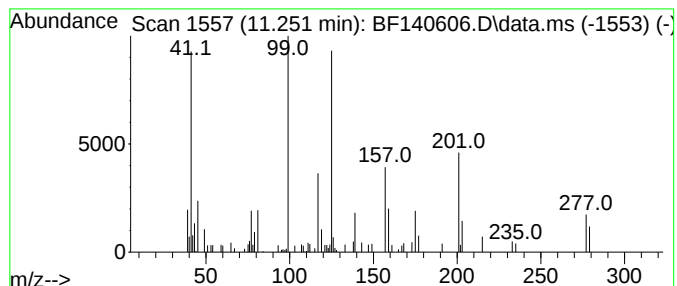
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Propanol, 1-chloro-, phos... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.251	2.41 ng	77569	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	72
2			Tris(2-chloropropyl) phosphate	326	C9H18Cl3O4P	006145-73-9	72
3			Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	52
4			Tris(3-chloropropyl) phosphate	326	C9H18Cl3O4P	001067-98-7	35
5			Bis(3-chloro-1-propyl)(1-chloro-...	326	C9H18Cl3O4P	137888-35-8	35



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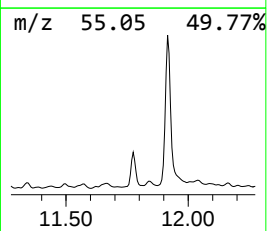
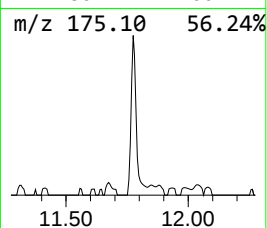
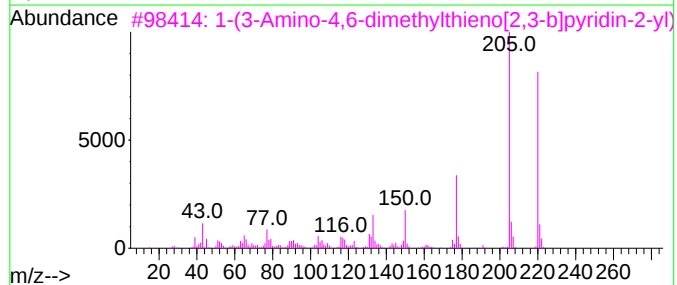
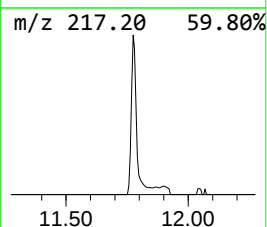
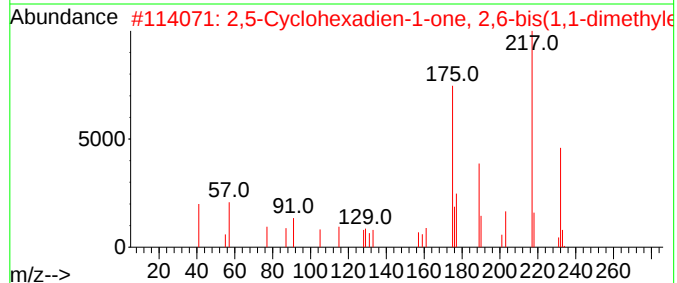
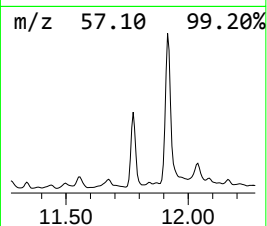
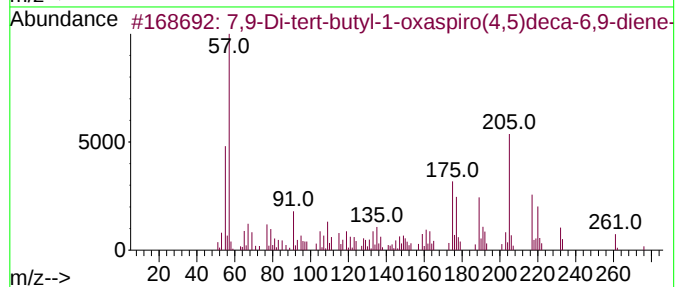
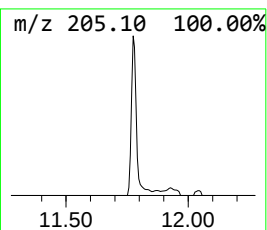
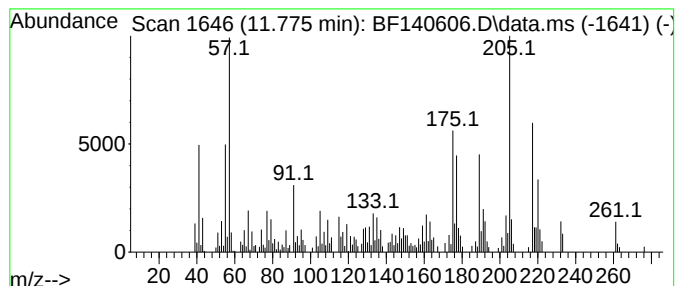
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.775	5.18 ng	166454	Phenanthrene-d10	11.392	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99
2	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	80
3	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	45
4	2,5-di-tert-Butyl-1,4-benzoquinone	220	C14H20O2	002460-77-7	44
5	1-Methyl-1H-pyrazolo[3,4-b]pyrid...	220	C10H16N4Si	1000468-34-6	42



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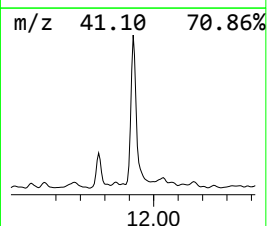
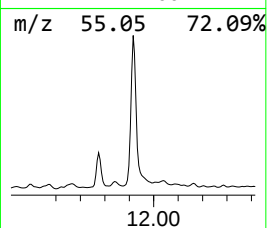
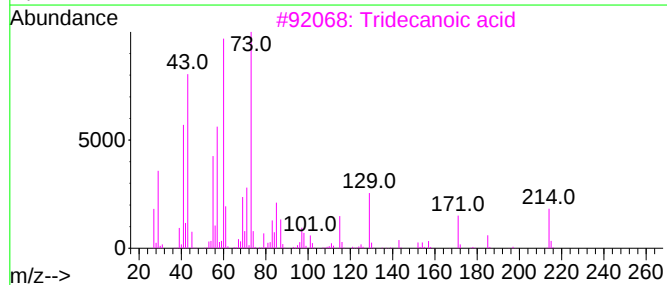
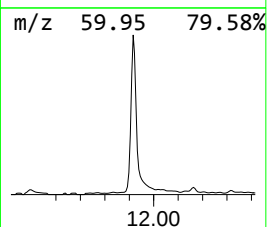
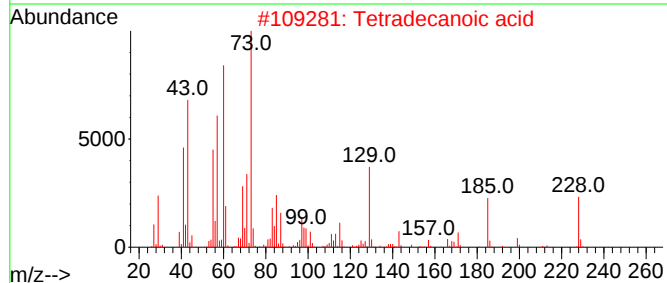
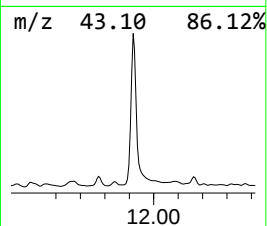
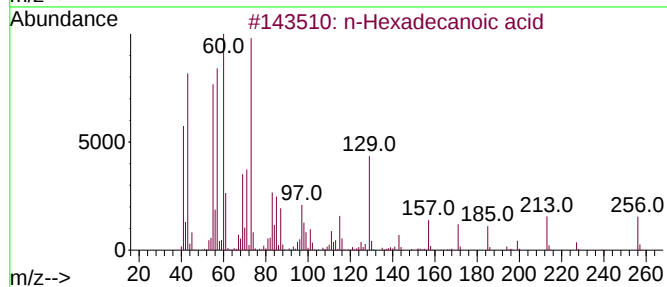
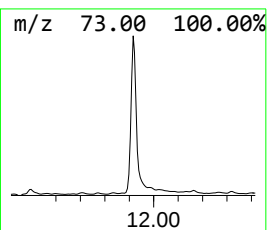
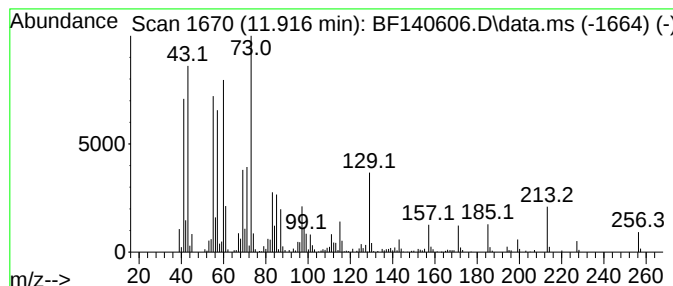
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.916	12.98 ng	417333	Phenanthrene-d10		11.392	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid		228	C14H28O2	000544-63-8	94
3	Tridecanoic acid		214	C13H26O2	000638-53-9	91
4	Pentadecanoic acid		242	C15H30O2	001002-84-2	91
5	Dodecanoic acid		200	C12H24O2	000143-07-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
Operator : RC/JU
Sample : P4947-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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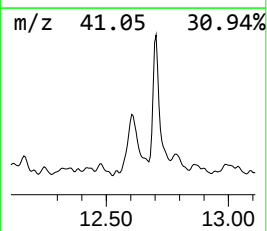
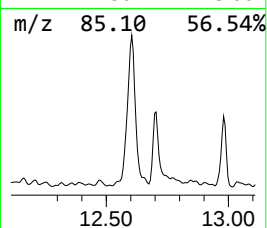
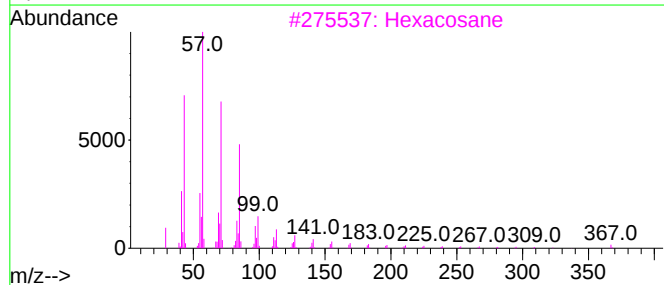
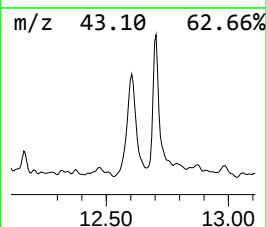
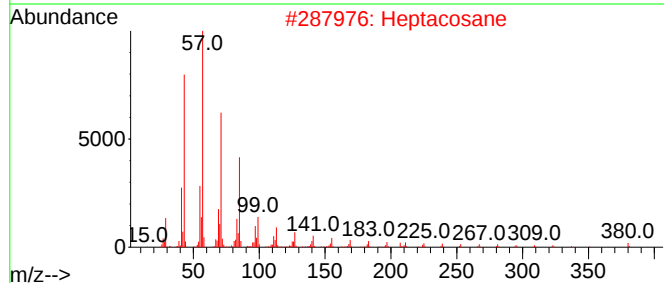
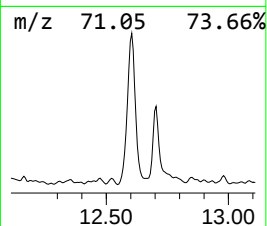
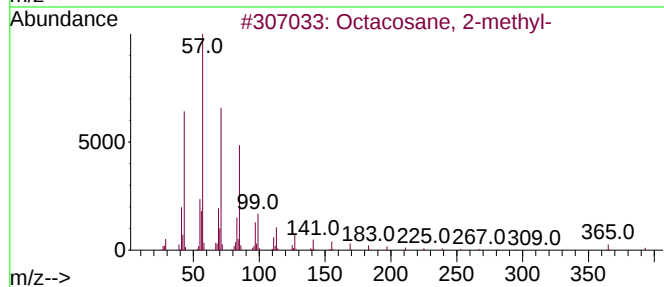
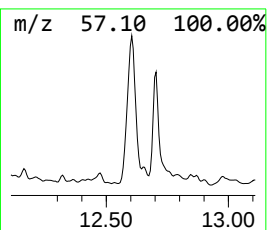
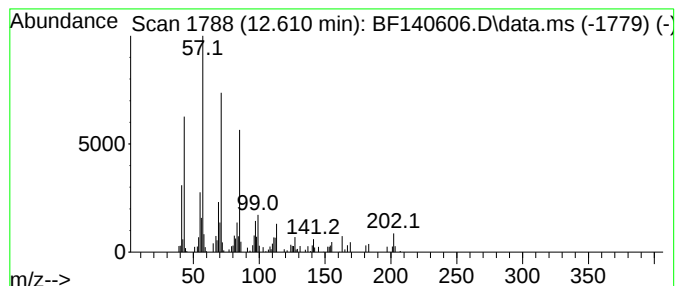
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octacosane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	4.61 ng	148055	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2			Heptacosane	380	C27H56	000593-49-7	90
3			Hexacosane	366	C26H54	000630-01-3	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
Operator : RC/JU
Sample : P4947-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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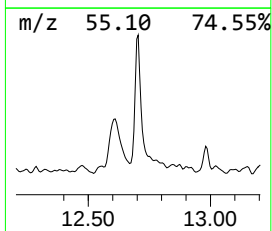
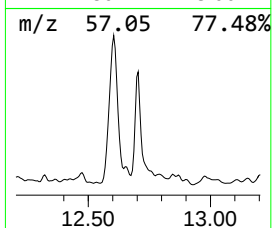
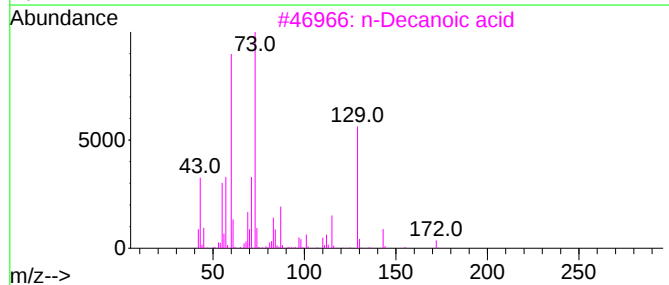
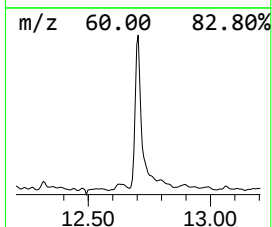
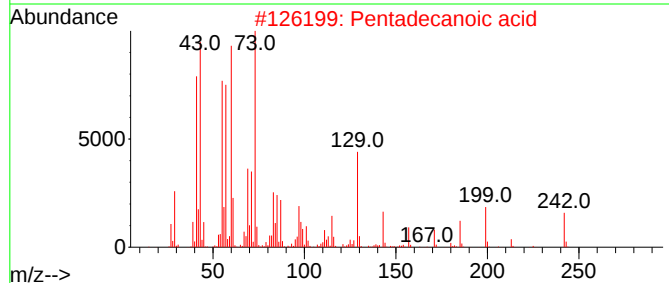
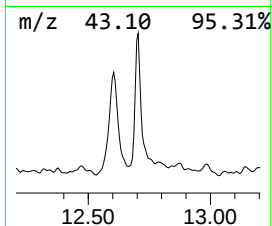
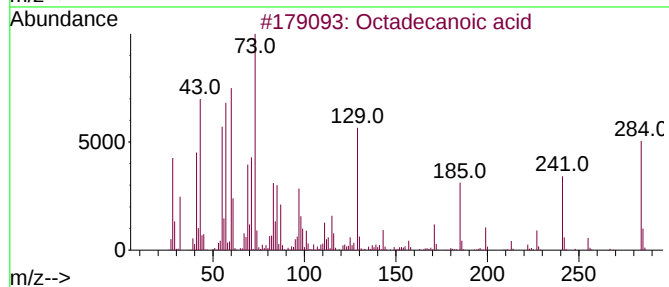
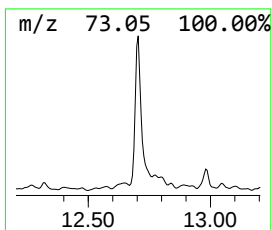
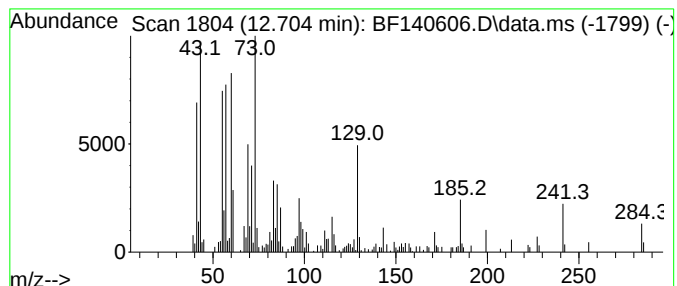
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.704	4.88 ng	156964	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3			n-Decanoic acid	172	C10H20O2	000334-48-5	86
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
Operator : RC/JU
Sample : P4947-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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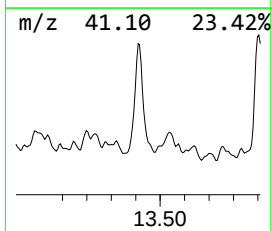
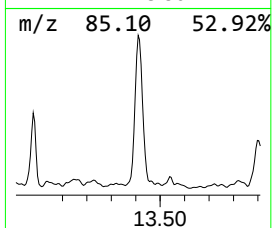
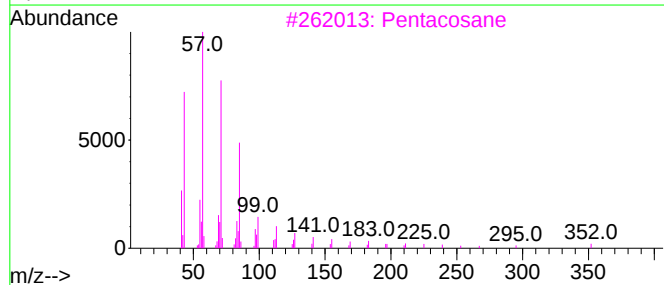
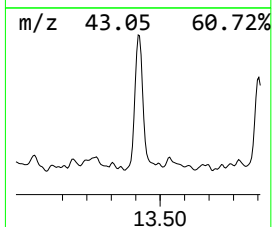
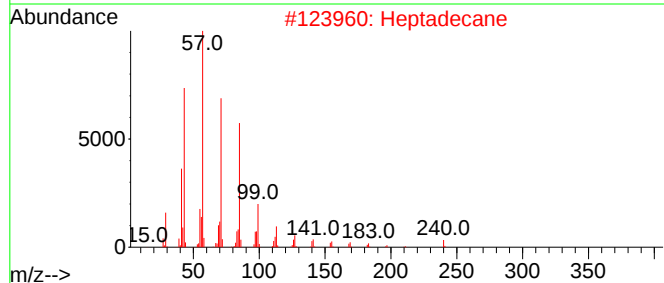
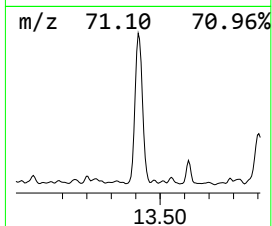
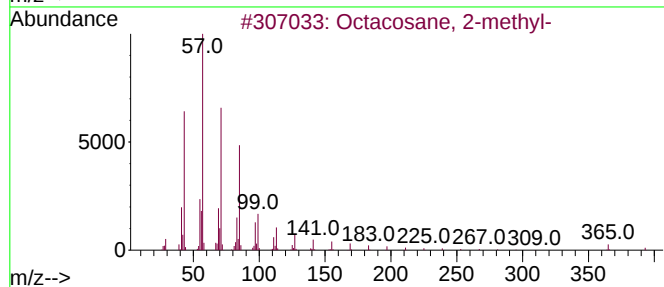
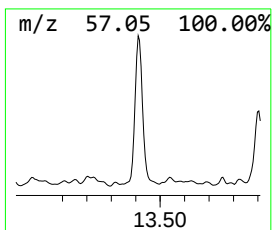
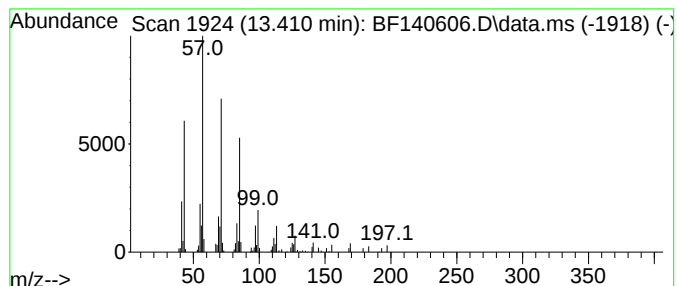
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.410	3.53 ng	100242	Chrysene-d12	14.045

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2		Heptadecane	240	C17H36	000629-78-7	91
3		Pentacosane	352	C25H52	000629-99-2	87
4		Hexacosane	366	C26H54	000630-01-3	87
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
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Sample : P4947-01
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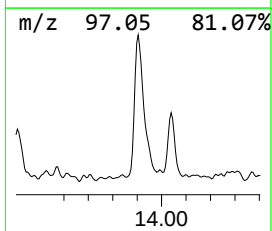
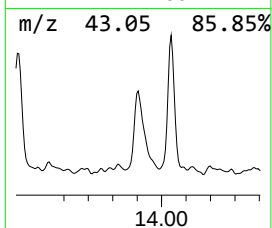
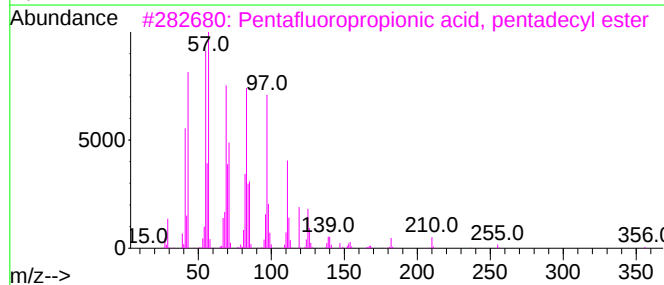
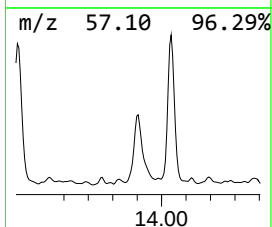
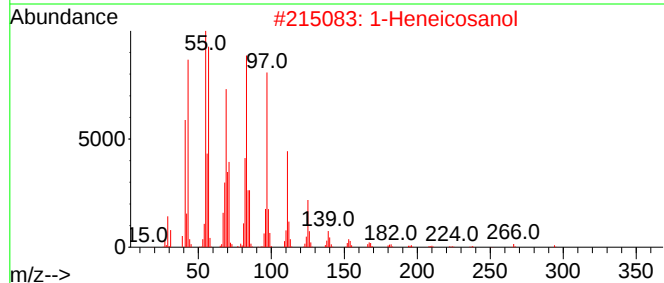
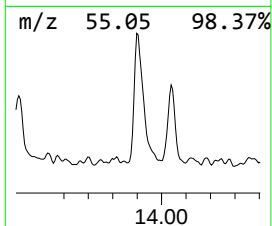
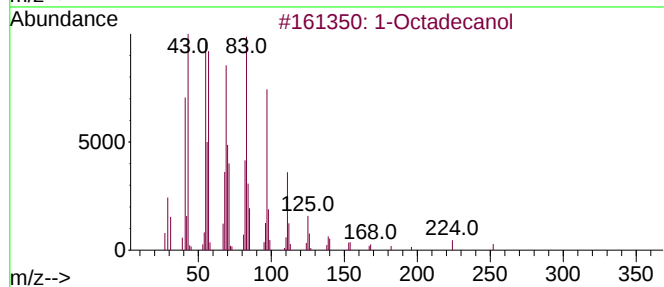
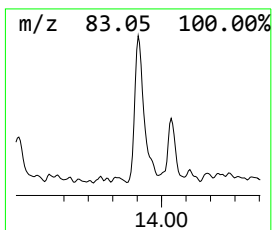
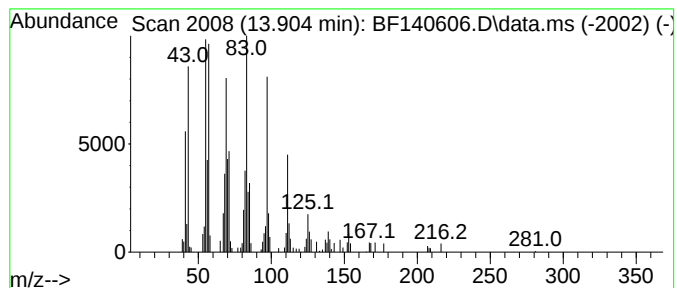
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Octadecanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.904	4.06 ng	115201	Chrysene-d12	14.045	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecanol	270	C18H38O	000112-92-5	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5	1-Hexacosene	364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
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Sample : P4947-01
Misc :
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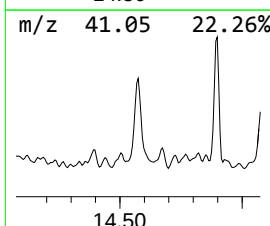
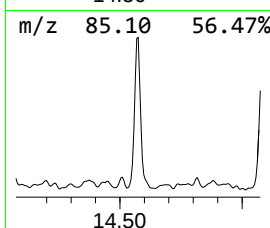
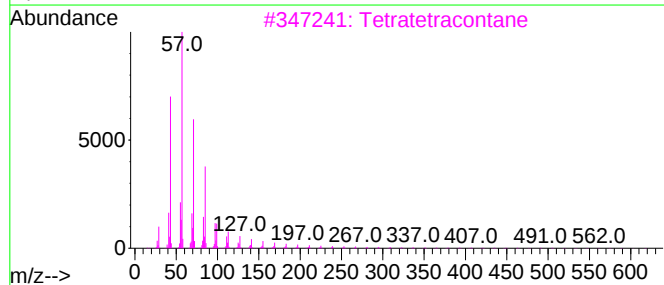
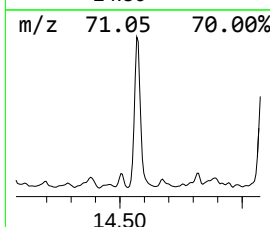
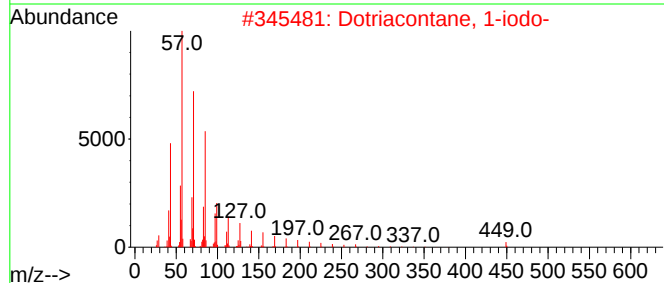
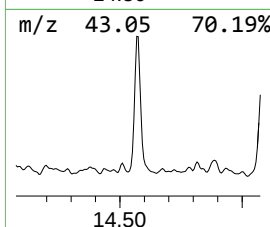
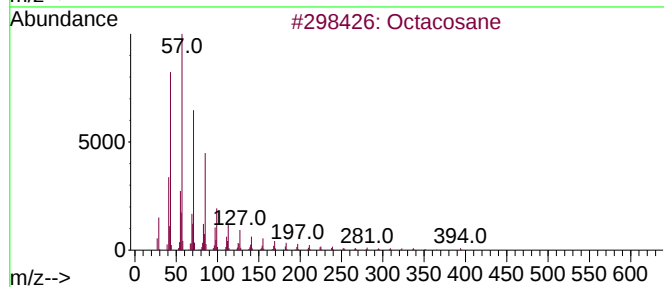
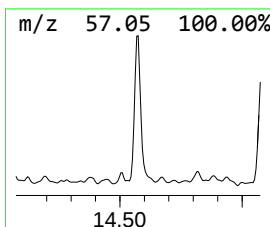
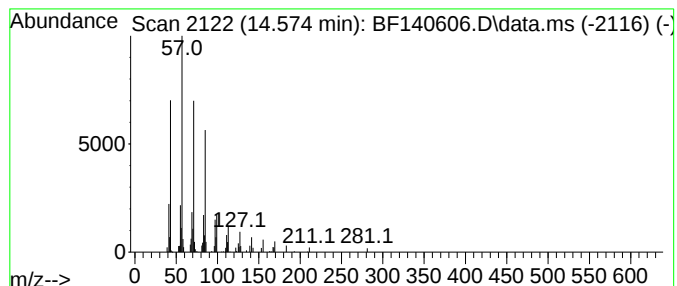
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.574	2.93 ng	83126	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3			Tetratetracontane	619	C44H90	007098-22-8	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
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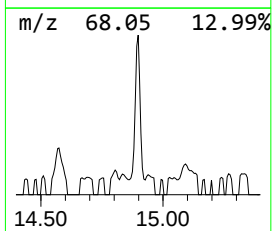
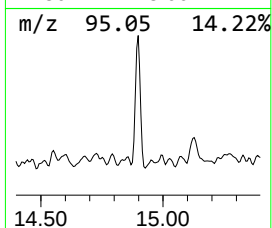
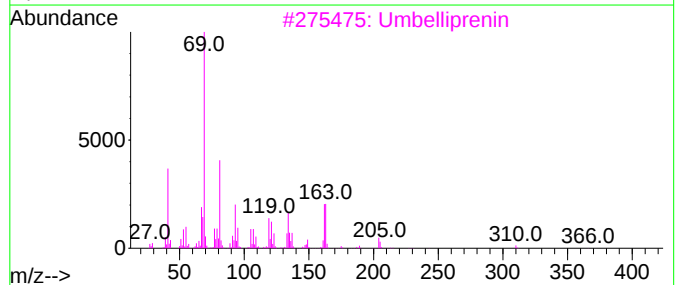
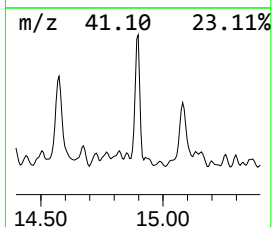
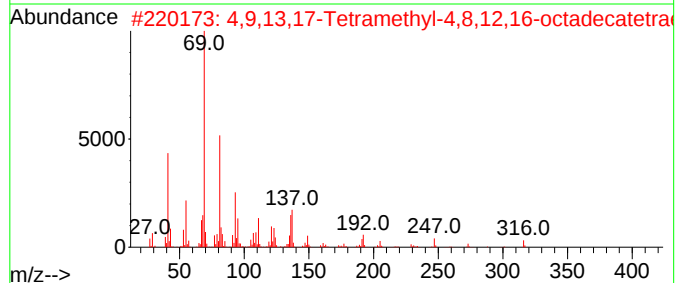
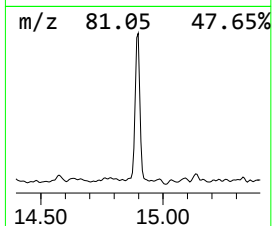
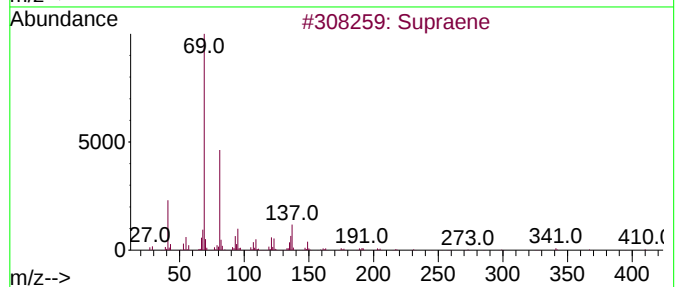
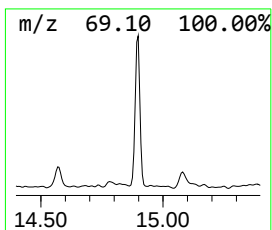
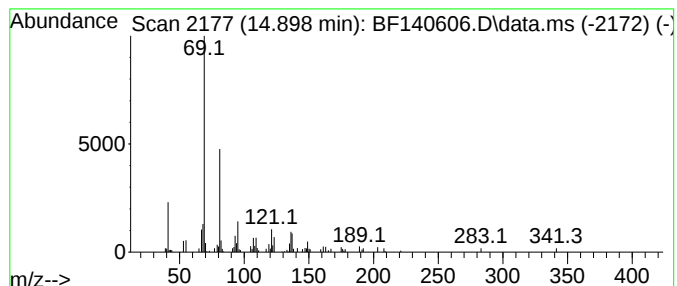
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Supraene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.898	3.24 ng	63491	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	86
2			4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	83
3			Umbelliprenin	366	C24H30O3	023838-17-7	72
4			4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64
5			3,7-Dimethyl-1-phenylsulfonyl-2,...	278	C16H22O2S	1000432-27-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
Operator : RC/JU
Sample : P4947-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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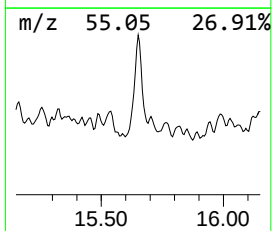
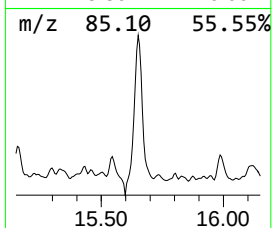
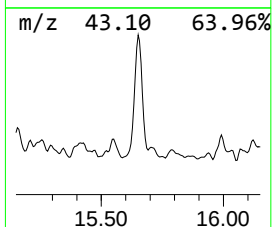
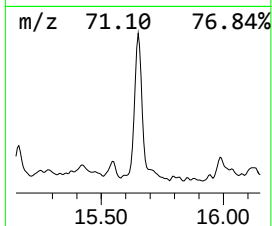
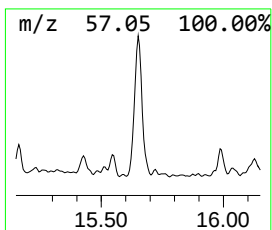
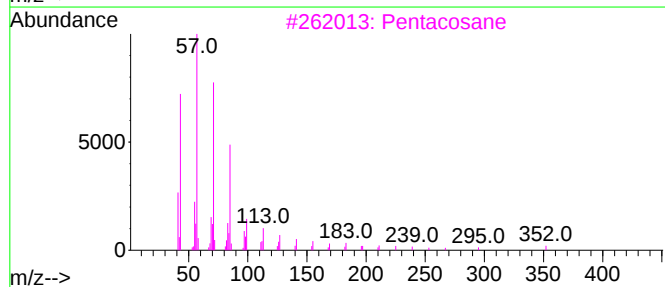
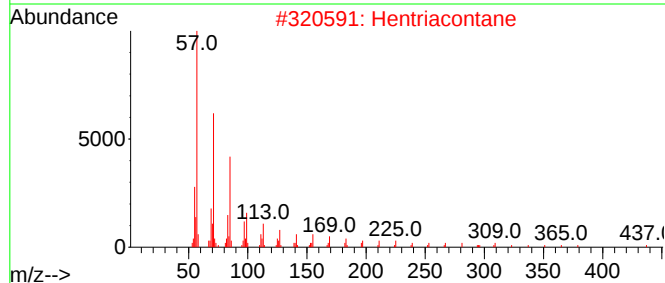
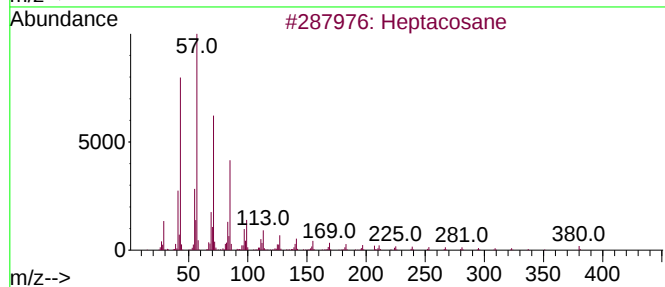
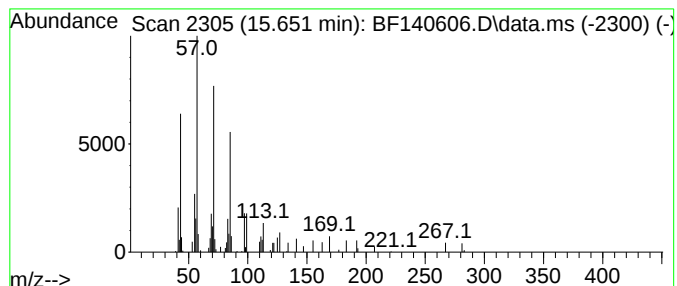
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.651	2.30 ng	45179	Perylene-d12	15.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	91
2		Hentriacontane	437	C31H64	000630-04-6	90
3		Pentacosane	352	C25H52	000629-99-2	90
4		Heneicosane	296	C21H44	000629-94-7	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
Operator : RC/JU
Sample : P4947-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
A3988

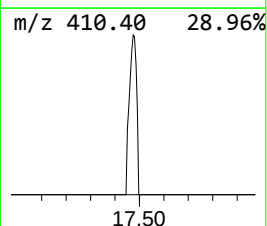
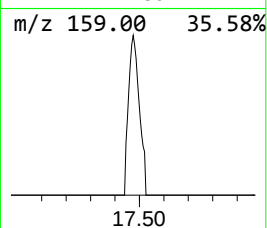
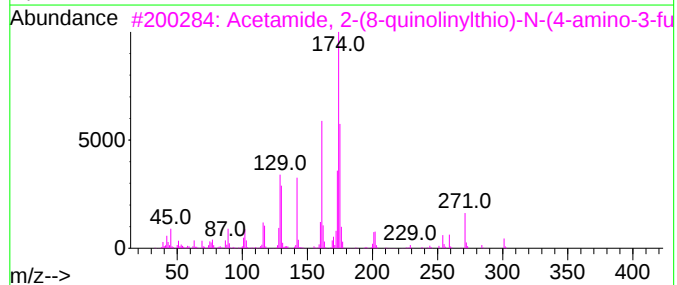
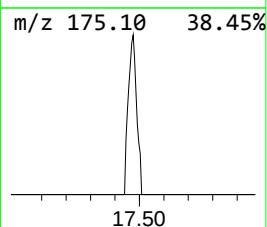
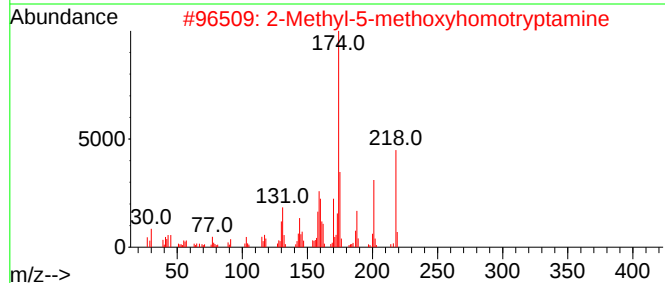
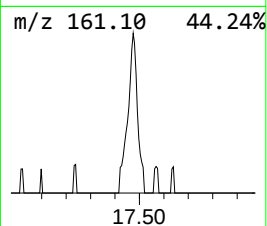
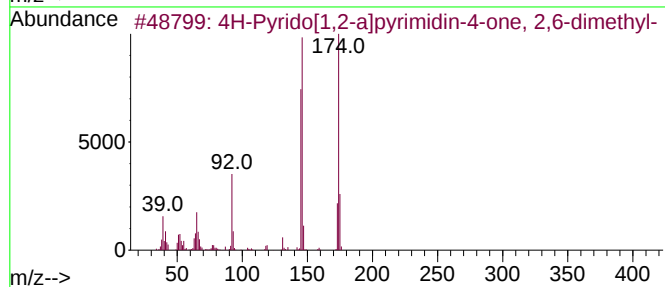
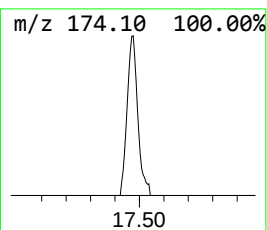
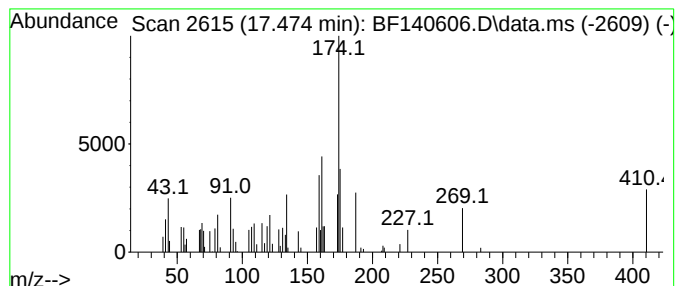
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown17.474 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.474	2.11 ng	41374	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Pyrido[1,2-a]pyrimidin-4-one,...	174	C10H10N2O	016867-28-0	38
2			2-Methyl-5-methoxyhomotryptamine	218	C13H18N2O	019017-58-4	37
3			Acetamide, 2-(8-quinolinylthio)-...	301	C13H11N5O2S	329780-80-5	35
4			Acetamide, N-[4-(chlorodifluorom...	394	C18H13ClF2N2O2S	1010309-89-6	32
5			Benzenemethanamine, 4-(trifluoro...	175	C8H8F3N	003300-51-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140606.D
Acq On : 25 Nov 2024 16:49
Operator : RC/JU
Sample : P4947-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
A3988

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	111.3	ng	2364340	1	6.869	424929	20.0
Benzophenone	10.616	8.8	ng	276390	3	9.904	625933	20.0
2-Propanol, 1-c...	11.251	2.4	ng	77569	4	11.392	642941	20.0
7,9-Di-tert-but...	11.775	5.2	ng	166454	4	11.392	642941	20.0
n-Hexadecanoic ...	11.916	13.0	ng	417333	4	11.392	642941	20.0
Octacosane, 2-m...	12.610	4.6	ng	148055	4	11.392	642941	20.0
Octadecanoic acid	12.704	4.9	ng	156964	4	11.392	642941	20.0
Heptadecane	13.410	3.5	ng	100242	5	14.045	567146	20.0
1-Octadecanol	13.904	4.1	ng	115201	5	14.045	567146	20.0
Octacosane	14.574	2.9	ng	83126	5	14.045	567146	20.0
Supraene	14.898	3.2	ng	63491	6	15.539	392435	20.0
Triacontane	15.080	3.0	ng	59931	6	15.539	392435	20.0
Heptacosane	15.651	2.3	ng	45179	6	15.539	392435	20.0
unknown17.474	17.474	2.1	ng	41374	6	15.539	392435	20.0