

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
 Data File : BF140932.D
 Acq On : 19 Dec 2024 17:49
 Operator : RC/JU
 Sample : P5343-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ281

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-BF121624.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF140932.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.493	573	578	606	rBV	701860	1123375	88.51%	9.660%
2	6.498	744	749	776	rBV	688124	1073819	84.61%	9.234%
3	6.845	803	808	817	rBV	267377	345862	27.25%	2.974%
4	7.416	899	905	927	rBV	433227	613784	48.36%	5.278%
5	7.675	945	949	964	rBV3	6032	14268	1.12%	0.123%
6	8.128	1021	1026	1042	rBV	325303	427354	33.67%	3.675%
7	9.198	1203	1208	1218	rBV	1006566	1269171	100.00%	10.914%
8	9.881	1319	1324	1329	rBV	336293	428891	33.79%	3.688%
9	10.592	1440	1445	1454	rVB	83114	108139	8.52%	0.930%
10	10.675	1454	1459	1472	rBV	452600	637133	50.20%	5.479%
11	11.375	1573	1578	1580	rBV	260812	346623	27.31%	2.981%
12	11.398	1580	1582	1587	rVV	146973	162358	12.79%	1.396%
13	11.451	1587	1591	1596	rVV	12421	16664	1.31%	0.143%
14	11.892	1659	1666	1674	rBV3	79254	181419	14.29%	1.560%
15	11.986	1678	1682	1691	rVB	39394	77467	6.10%	0.666%
16	12.169	1709	1713	1717	rBV	13281	15972	1.26%	0.137%
17	12.210	1717	1720	1725	rVB2	22120	28523	2.25%	0.245%
18	12.427	1752	1757	1760	rBV	22578	32502	2.56%	0.280%
19	12.522	1769	1773	1780	rVB3	17166	30029	2.37%	0.258%
20	12.592	1780	1785	1790	rBV	232329	303851	23.94%	2.613%
21	12.680	1794	1800	1806	rVV6	16211	38294	3.02%	0.329%
22	12.739	1806	1810	1814	rVV2	19013	28867	2.27%	0.248%
23	12.822	1817	1824	1829	rVV	218382	323992	25.53%	2.786%
24	12.874	1829	1833	1839	rVV3	13161	21980	1.73%	0.189%
25	12.951	1841	1846	1852	rVV	689391	917763	72.31%	7.892%
26	13.039	1856	1861	1866	rBV3	19543	35568	2.80%	0.306%
27	13.151	1872	1880	1886	rBV	30340	62083	4.89%	0.534%
28	13.216	1886	1891	1894	rBV	15830	23410	1.84%	0.201%
29	13.257	1894	1898	1904	rVB2	25498	40344	3.18%	0.347%
30	13.351	1910	1914	1916	rBV	17417	23146	1.82%	0.199%
31	13.380	1916	1919	1929	rVB3	19400	32807	2.58%	0.282%
32	13.669	1965	1968	1974	rVB	23385	32505	2.56%	0.280%
33	13.780	1981	1987	1993	rBV2	36419	84269	6.64%	0.725%
34	13.851	1993	1999	2003	rVV3	60130	120830	9.52%	1.039%
35	13.886	2003	2005	2011	rVB2	28949	39623	3.12%	0.341%
36	14.033	2022	2030	2042	rVB2	332387	695474	54.80%	5.981%

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Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.433	2093	2098	2102	rBV	21244	35781	2.82%	0.308%
38	14.510	2102	2111	2118	rBV6	72999	192223	15.15%	1.653%
39	14.633	2129	2132	2136	rVB5	11343	16354	1.29%	0.141%
40	14.804	2159	2161	2165	rVB	20171	24425	1.92%	0.210%
41	14.951	2183	2186	2190	rVB3	20834	26794	2.11%	0.230%
42	14.998	2191	2194	2198	rVB	18845	24124	1.90%	0.207%
43	15.110	2207	2213	2216	rBV2	120121	220923	17.41%	1.900%
44	15.180	2221	2225	2242	rVB4	109689	289032	22.77%	2.486%
45	15.415	2261	2265	2271	rBV	63373	93919	7.40%	0.808%
46	15.480	2272	2276	2281	rVV	85793	135114	10.65%	1.162%
47	15.545	2282	2287	2300	rVB	220990	377818	29.77%	3.249%
48	15.957	2353	2357	2363	rBV	30289	48974	3.86%	0.421%
49	16.033	2364	2370	2382	rBV8	23984	87684	6.91%	0.754%
50	16.221	2396	2402	2417	rVB8	21031	71085	5.60%	0.611%
51	16.421	2431	2436	2440	rBV7	22653	38941	3.07%	0.335%
52	17.062	2539	2545	2553	rVB	54047	121450	9.57%	1.044%
53	17.521	2618	2623	2629	rBV	29905	65747	5.18%	0.565%

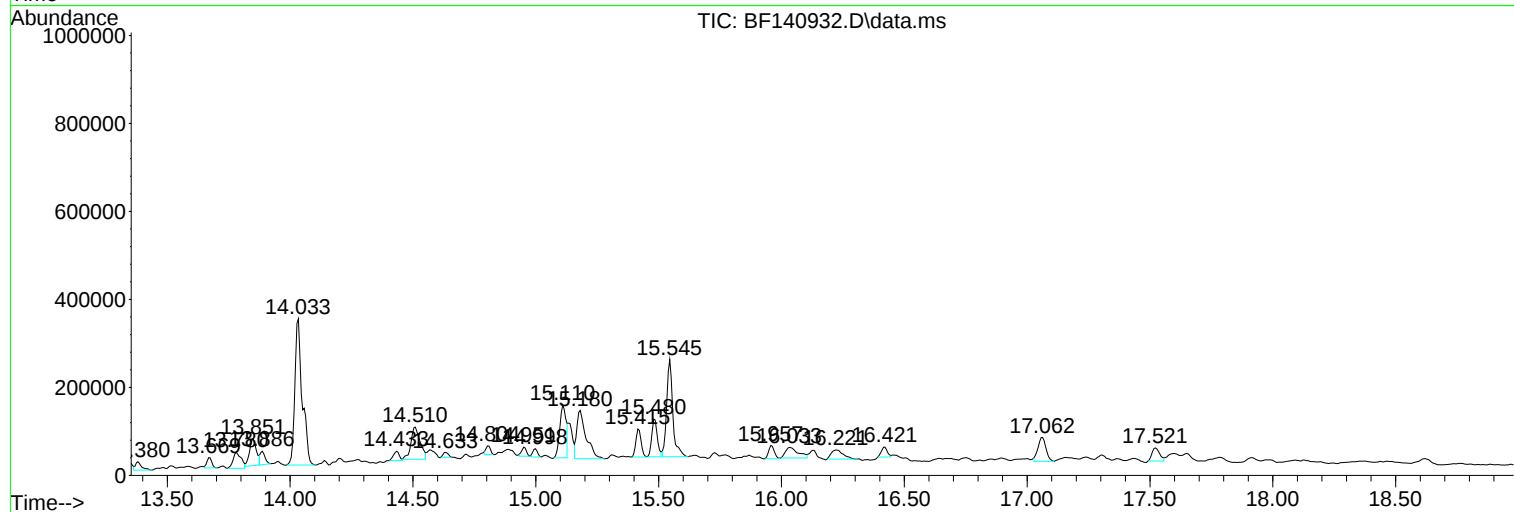
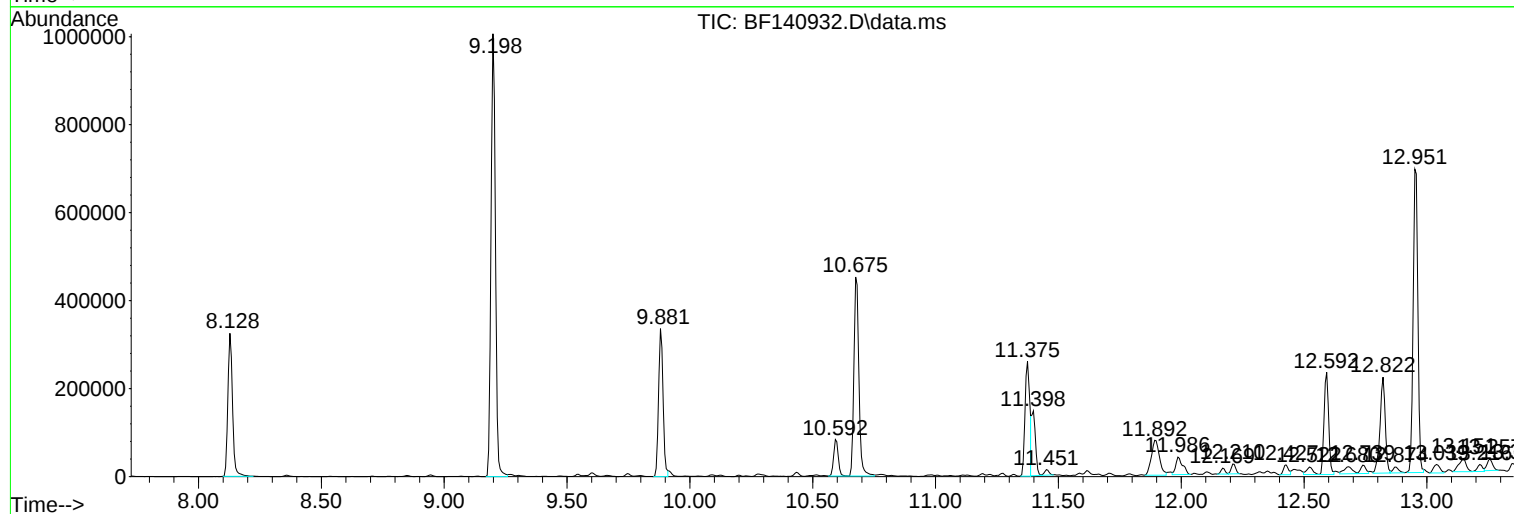
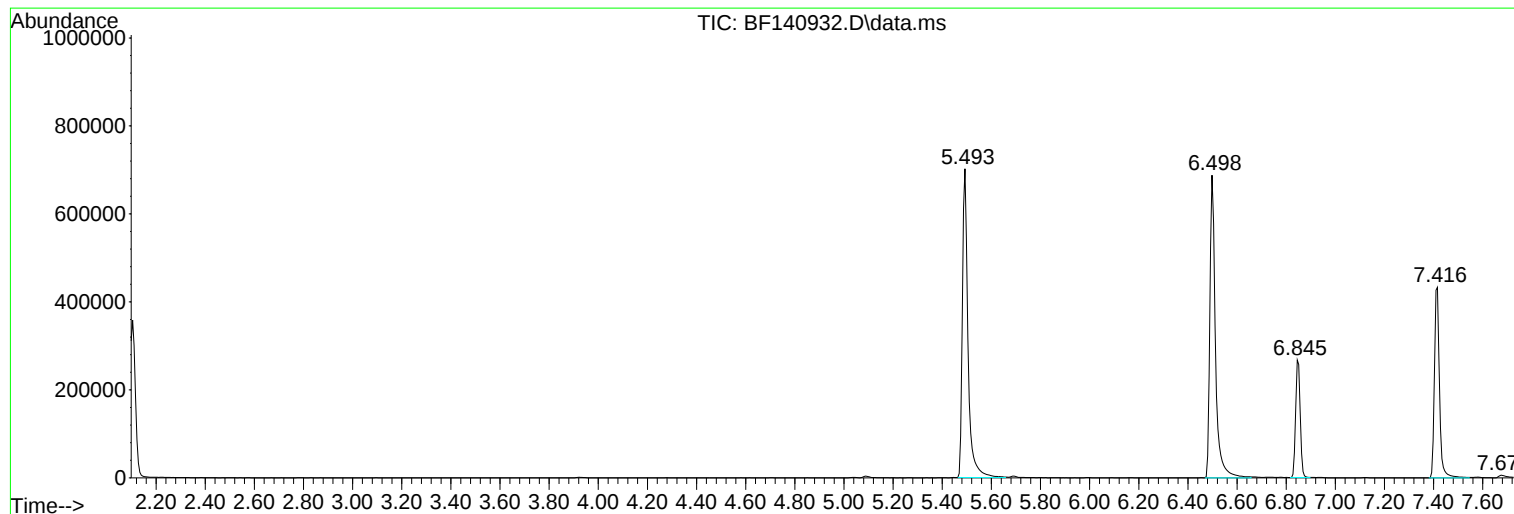
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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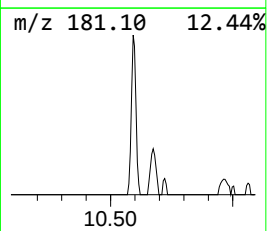
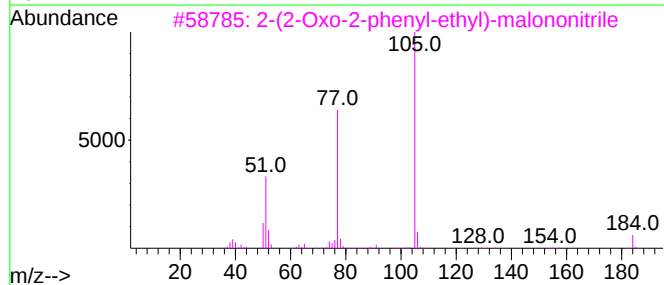
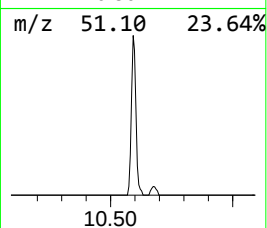
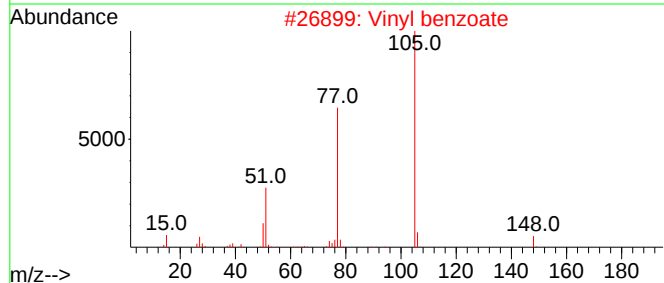
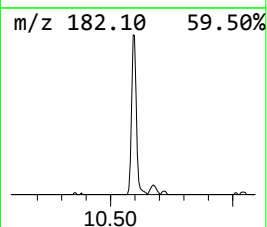
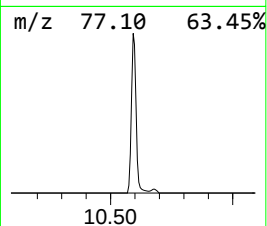
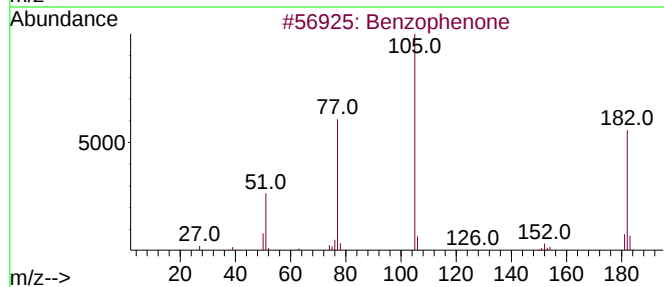
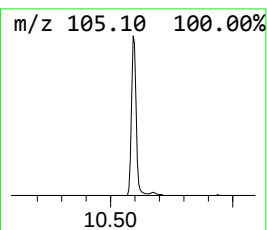
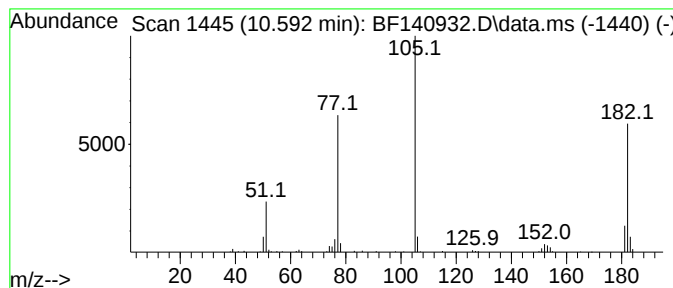
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.592	5.04 ng	108139	Acenaphthene-d10	9.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Vinyl benzoate	148	C9H8O2	000769-78-8	47
3			2-(2-Oxo-2-phenyl-ethyl)-malononitrile	184	C11H8N2O	1000296-76-9	47
4			Benzoyl isothiocyanate	163	C8H5NOS	000532-55-8	47
5			Benzenecarbothioic acid, S-methyl...	152	C8H8OS	005925-68-8	47



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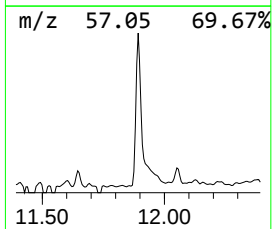
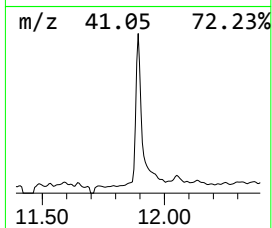
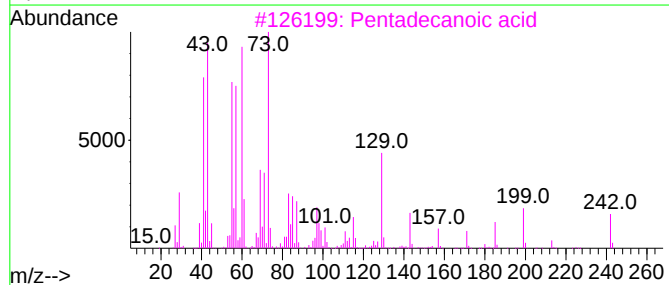
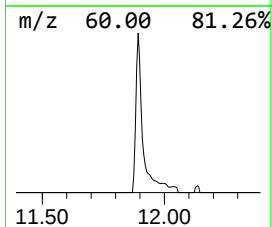
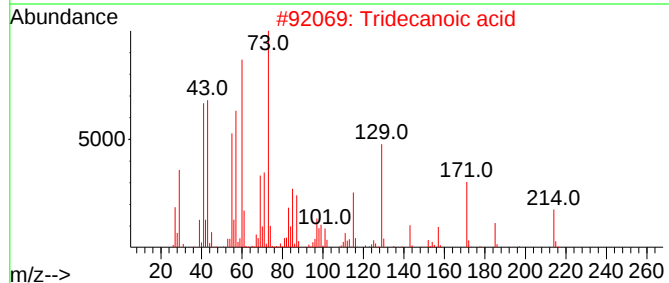
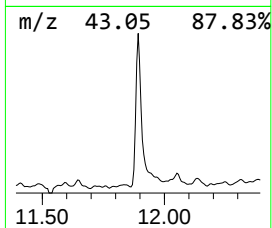
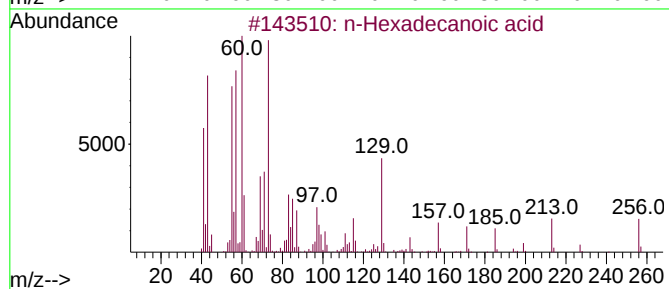
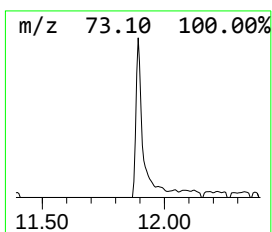
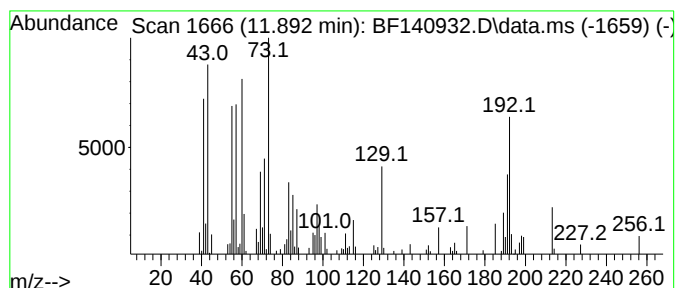
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.892	10.47 ng	181419	Phenanthrene-d10	11.374	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	70
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	60
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	25
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	25



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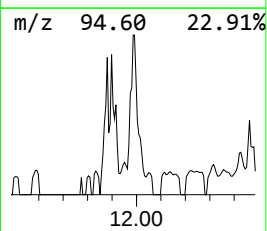
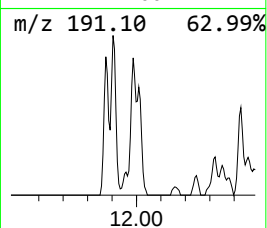
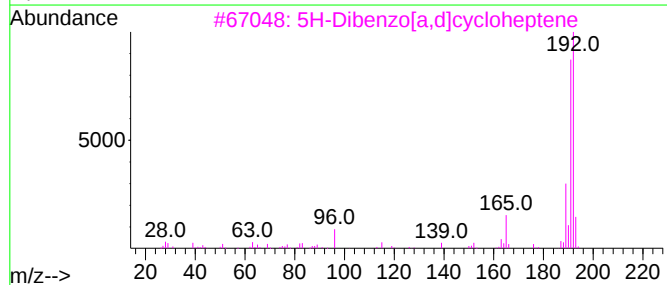
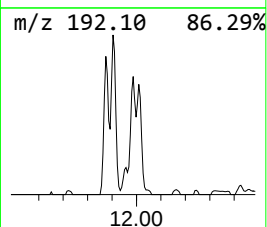
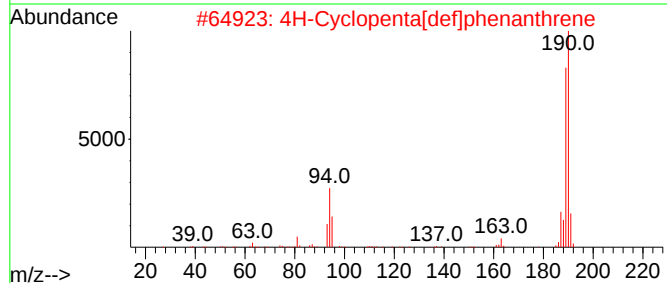
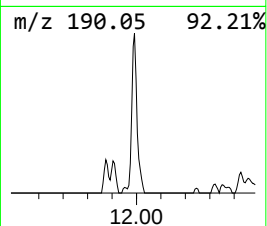
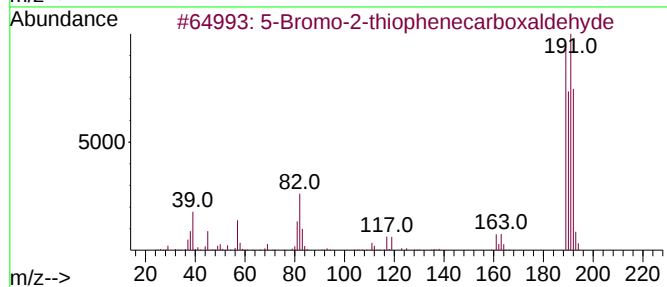
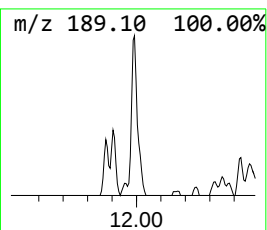
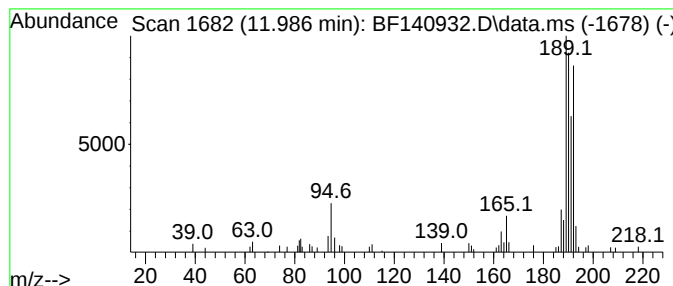
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 5-Bromo-2-thiophenecarboxal... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.986	4.47 ng	77467	Phenanthrene-d10	11.374	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	58
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
3	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	41
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38



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Misc :
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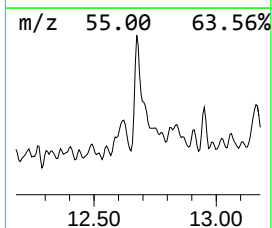
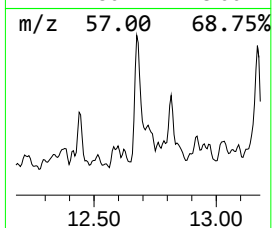
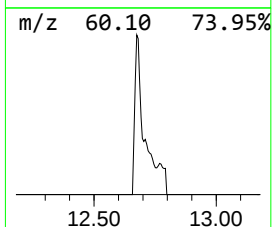
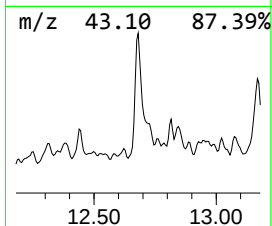
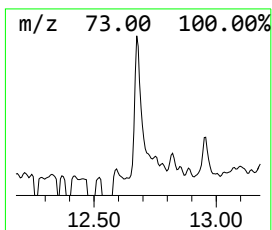
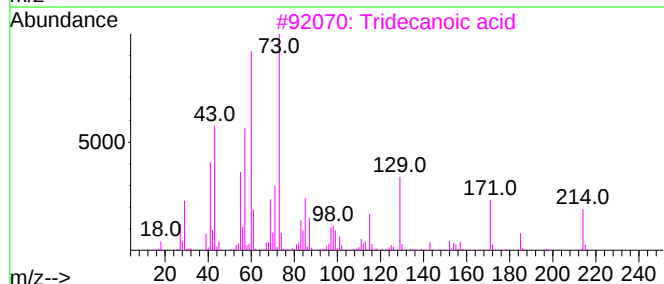
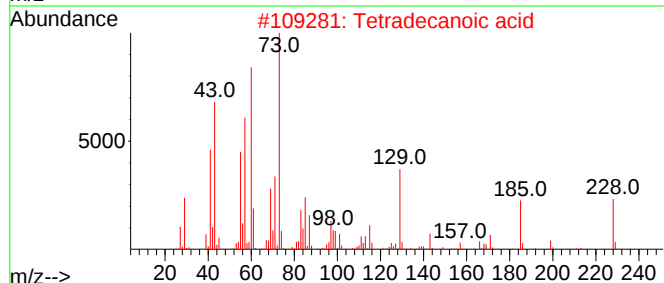
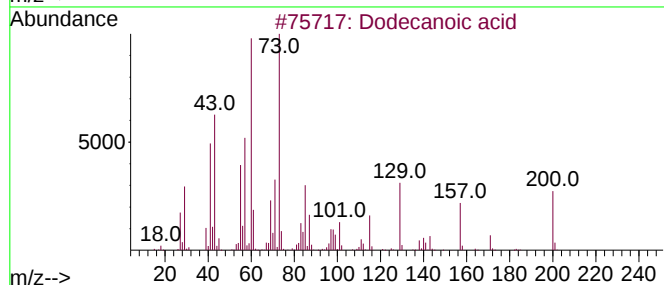
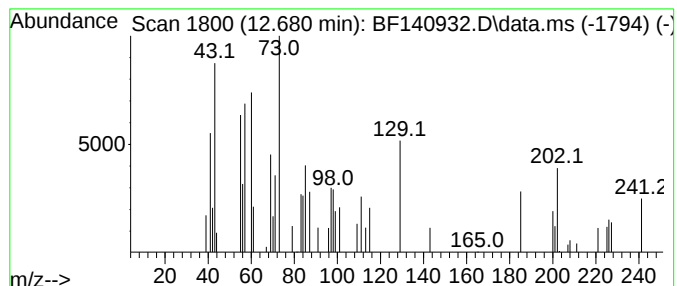
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 unknown12.680 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.680	2.21 ng	38294	Phenanthrene-d10	11.374	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecanoic acid	200	C12H24O2	000143-07-7	47
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	43
3	Tridecanoic acid	214	C13H26O2	000638-53-9	38
4	.alpha.-d-Riboside, 1-O-dodecyl-	318	C17H34O5	1000154-76-8	27
5	8-Methylnonanoic acid	172	C10H20O2	005963-14-4	27



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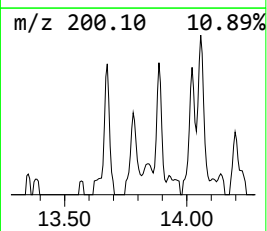
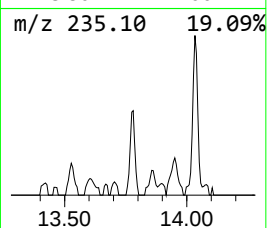
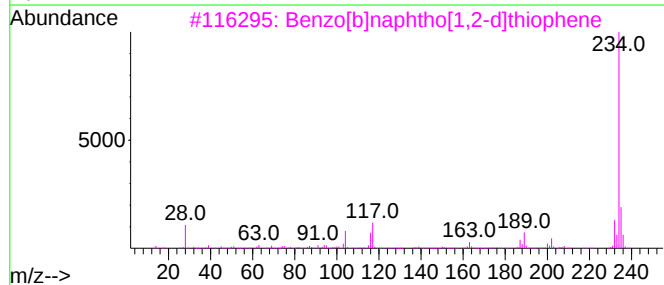
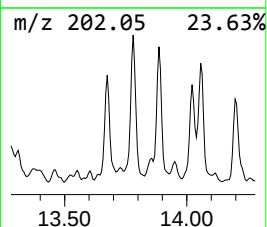
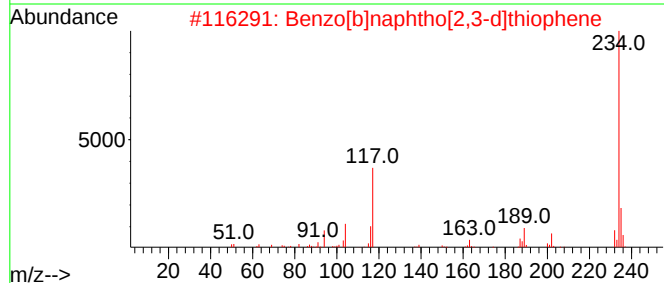
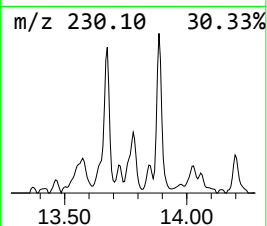
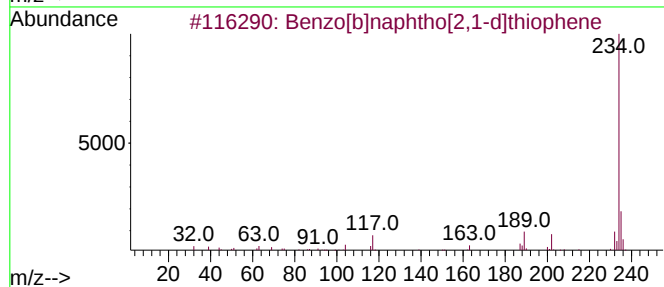
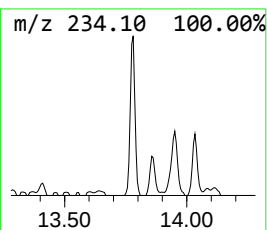
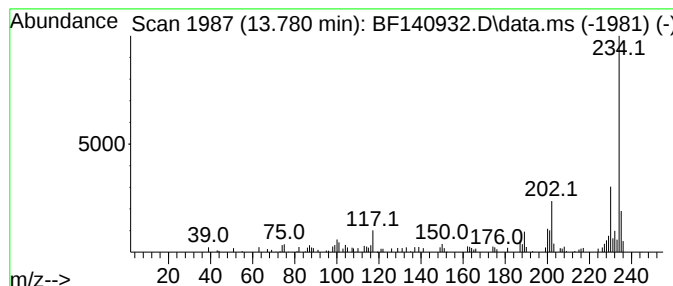
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	2.42 ng	84269	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
2			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	62
3			Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	58
4			Quinoxalino[2,3-b]quinoxaline, 5...	234	C14H10N4	000531-46-4	47
5			pyrido[1',2':1,2]imidazo[4,5-b]q...	234	C14H10N4	1000401-06-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ281

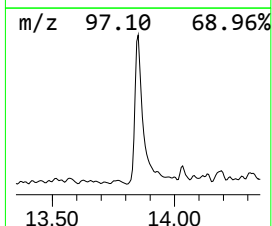
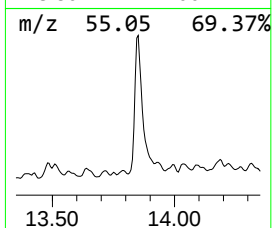
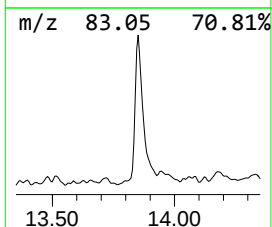
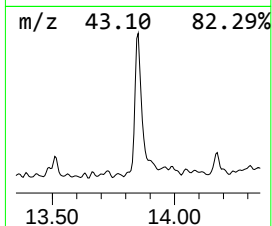
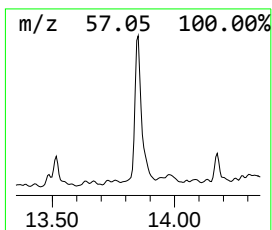
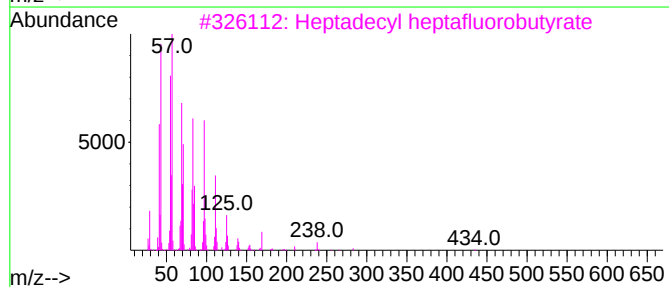
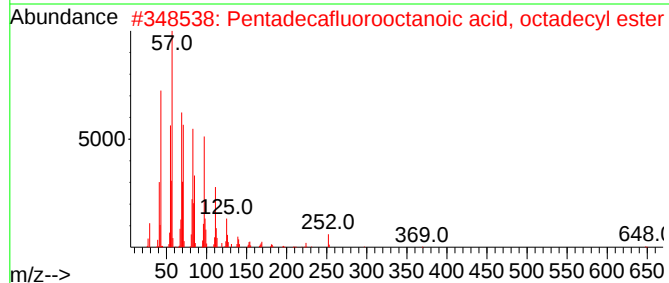
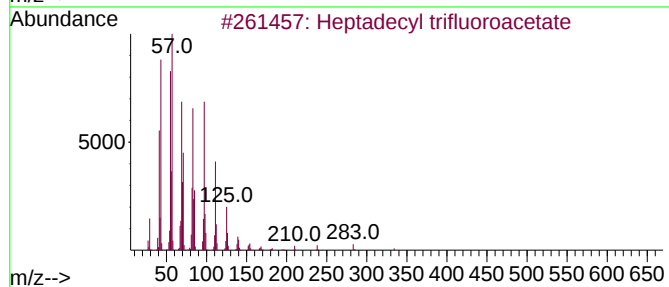
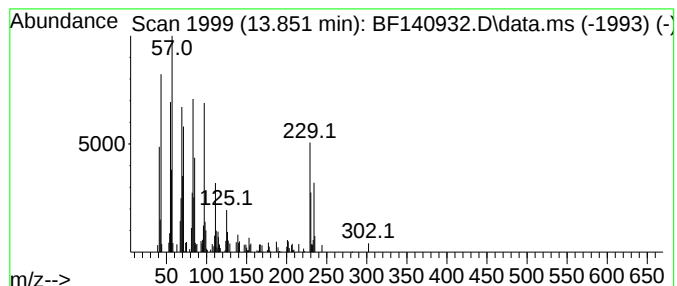
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptadecyl trifluoroacetate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.851	3.47 ng	120830	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	90
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	83
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	78
4			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	64
5			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

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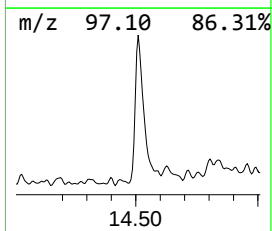
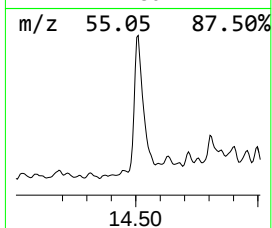
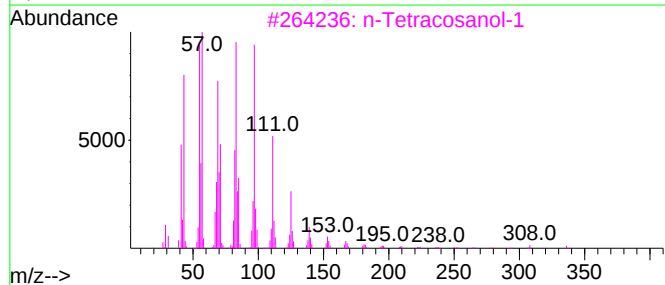
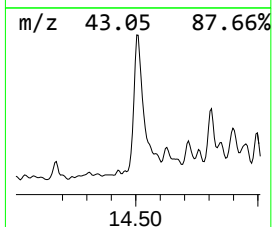
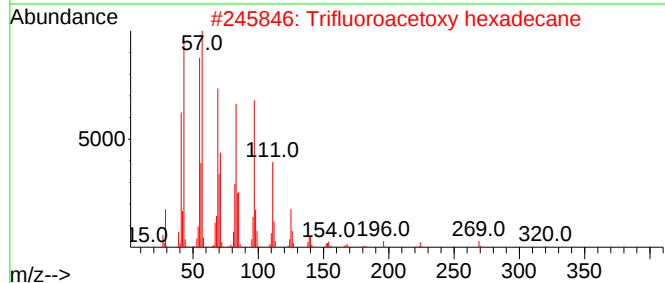
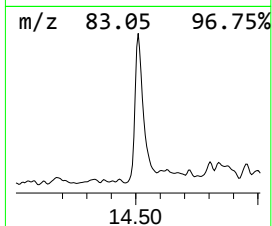
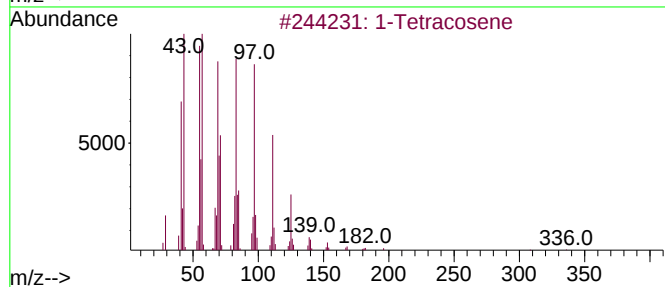
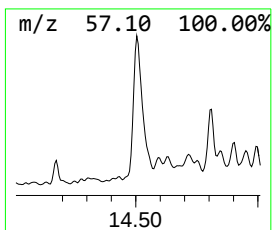
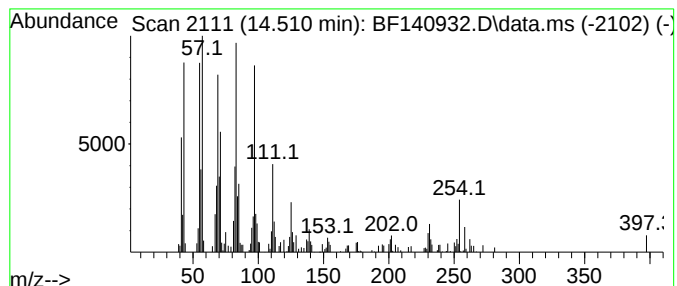
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Tetracosene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.510	5.53 ng	192223	Chrysene-d12	14.033

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene		336	C24H48	010192-32-2	93
2	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	93
3	n-Tetracosanol-1		354	C24H50O	000506-51-4	93
4	1-Heneicosanol		312	C21H44O	015594-90-8	93
5	1-Hexacosene		364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

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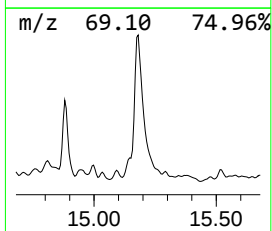
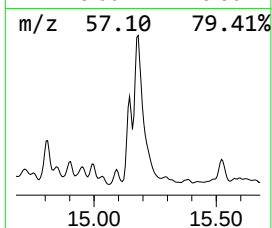
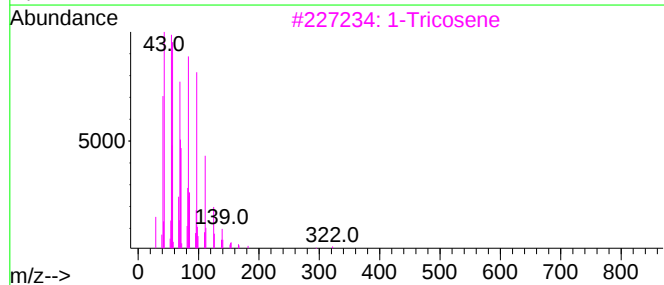
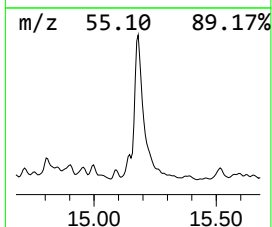
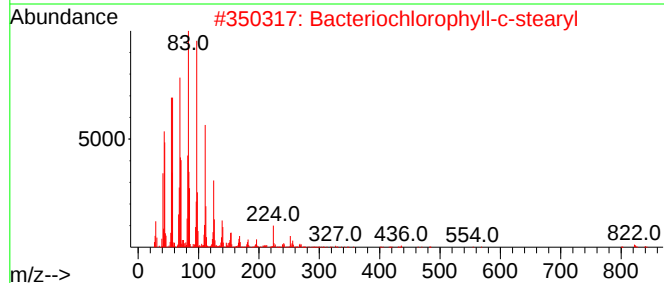
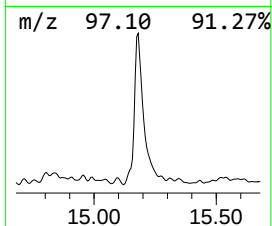
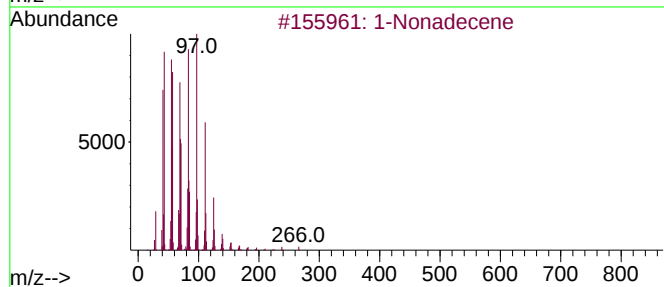
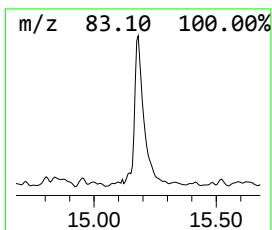
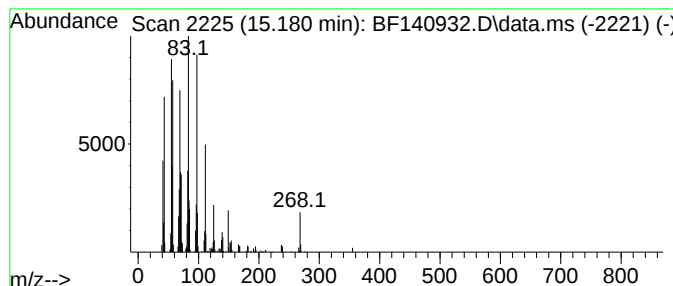
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Nonadecene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.180	15.30 ng	289032	Perylene-d12	15.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene			266	C19H38	018435-45-5	90
2	Bacteriochlorophyll-c-stearyl			841	C52H72MgN4O4	1010164-49-7	87
3	1-Tricosene			322	C23H46	018835-32-0	74
4	1-Tetracosene			336	C24H48	010192-32-2	70
5	1-Octadecanol			270	C18H38O	000112-92-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
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Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

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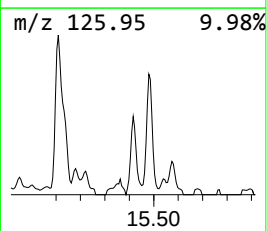
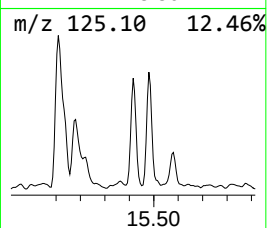
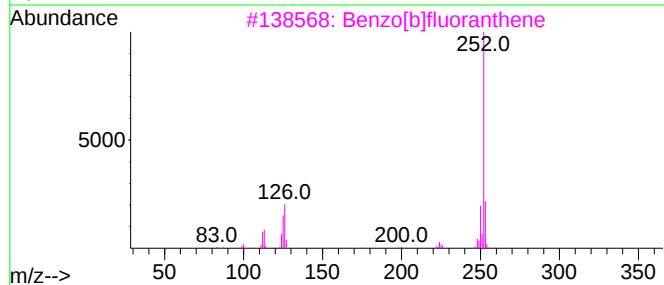
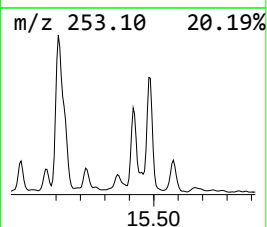
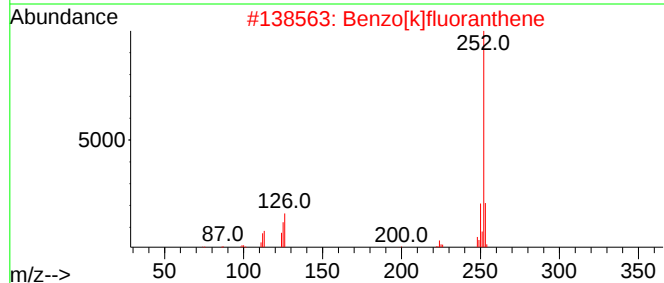
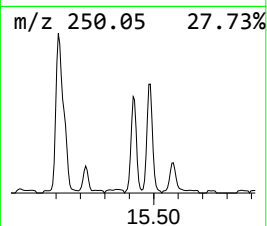
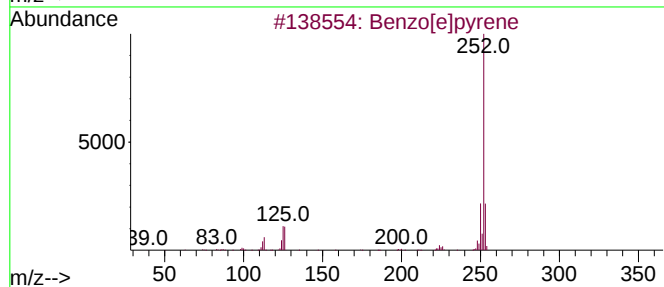
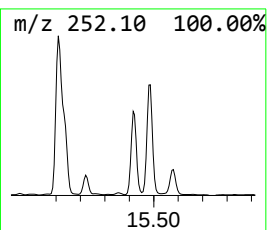
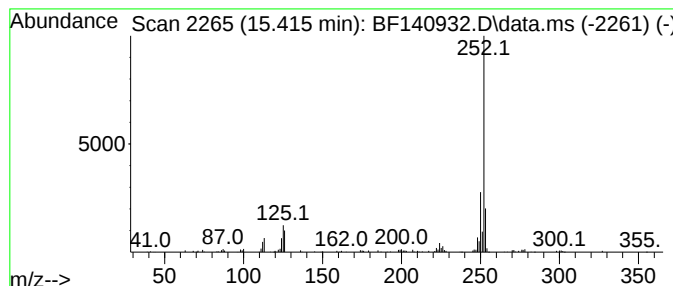
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.416	4.97 ng	93919	Perylene-d12	15.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
5			Perylene	252	C20H12	000198-55-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

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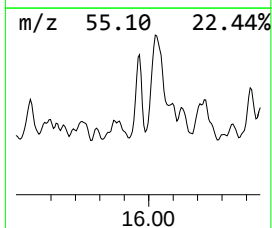
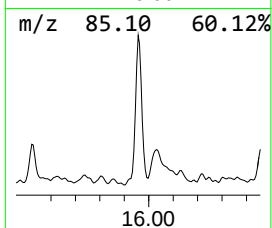
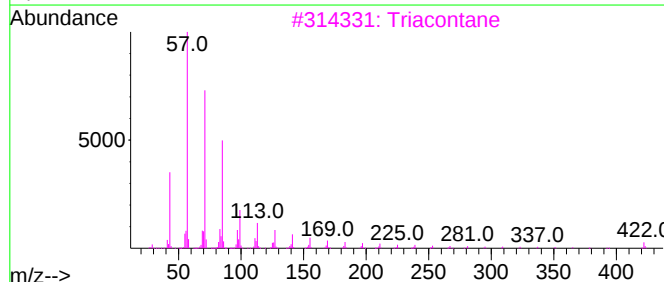
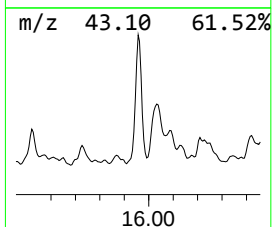
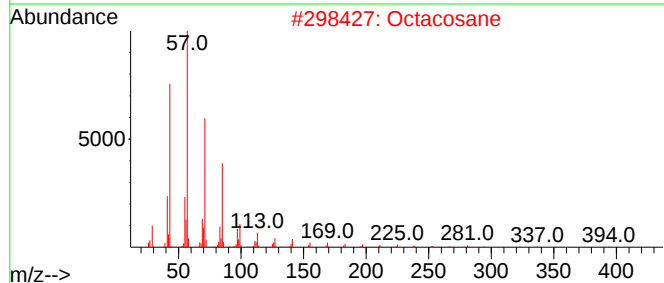
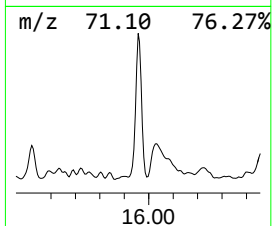
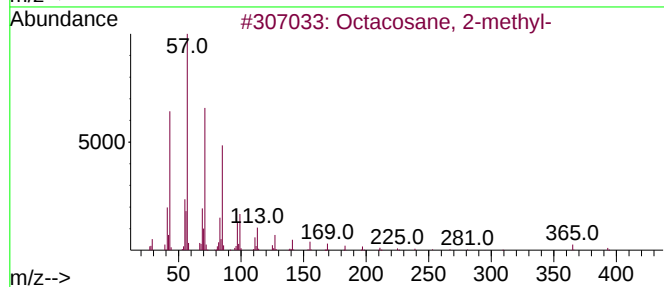
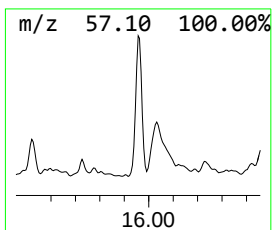
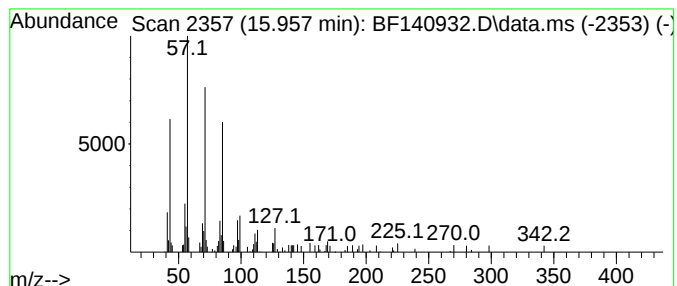
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octacosane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	2.59 ng	48974	Perylene-d12	15.545

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

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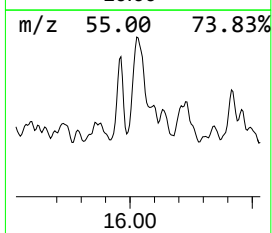
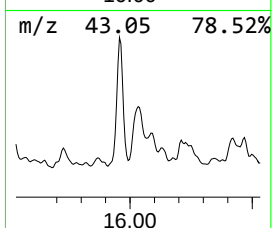
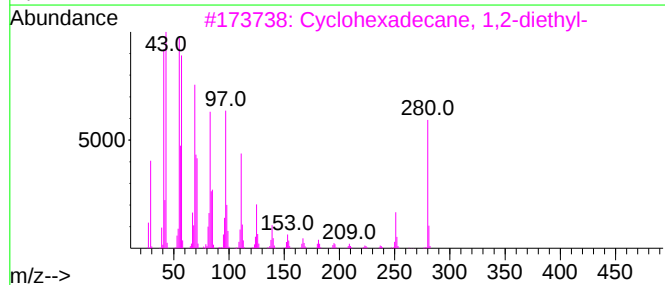
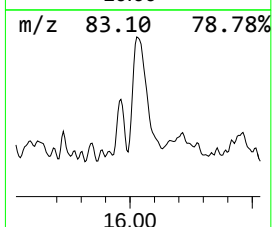
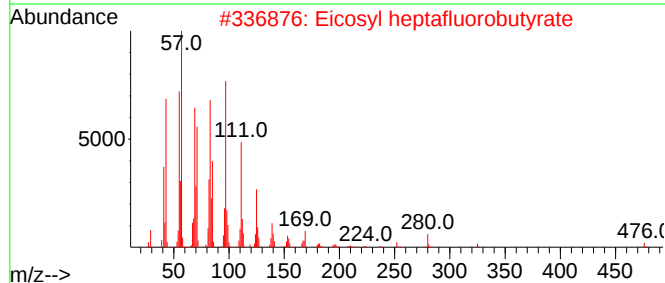
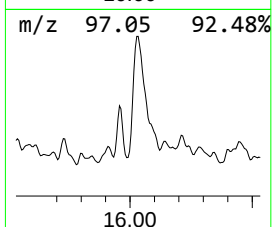
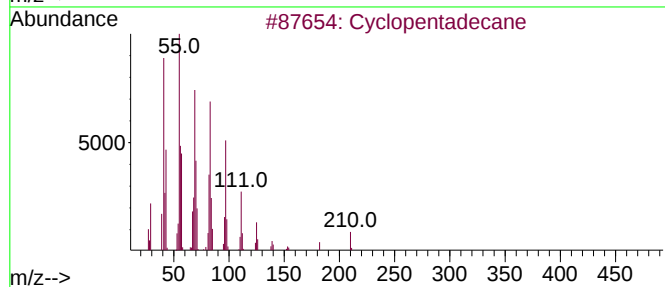
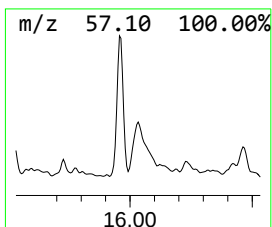
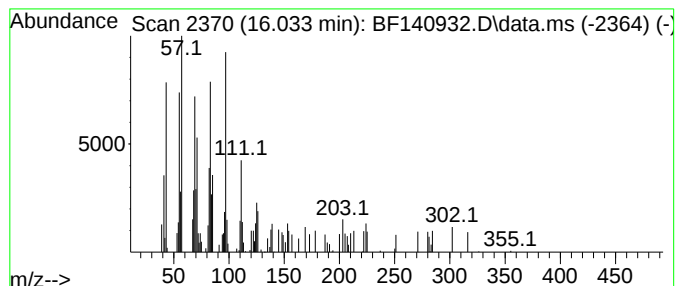
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclopentadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.033	4.64 ng	87684	Perylene-d12	15.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	92
2			Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	90
3			Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	89
4			1-Heptacosanol	396	C27H56O	002004-39-9	89
5			Octacosyl acetate	452	C30H60O2	018206-97-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ281

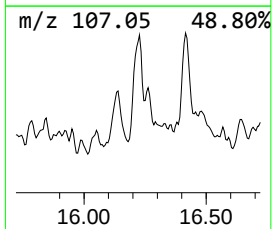
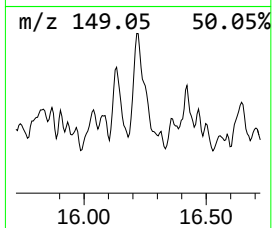
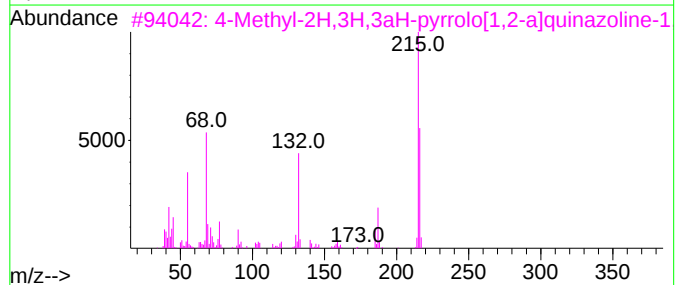
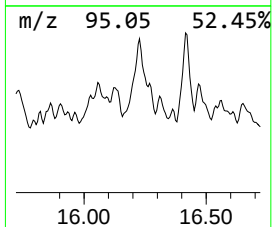
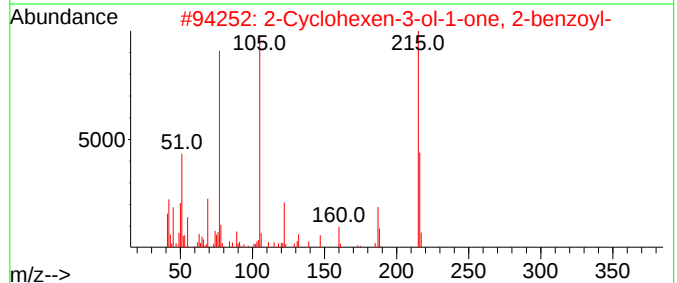
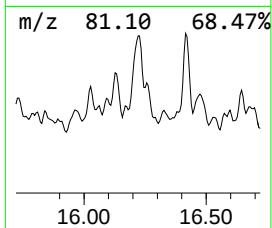
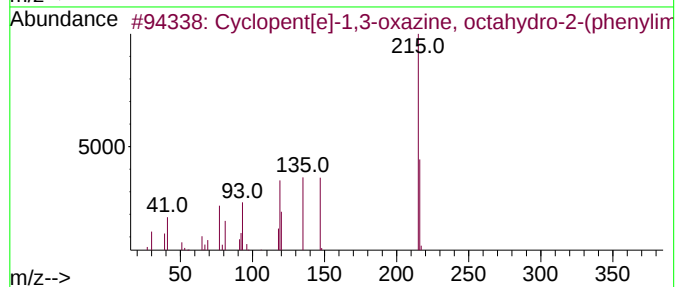
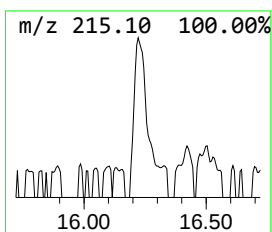
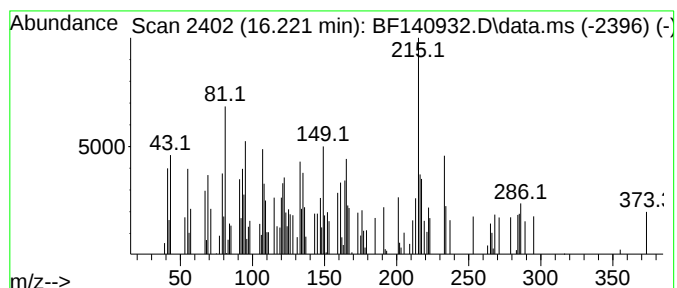
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown16.221 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.221	3.76 ng	71085	Perylene-d12	15.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopent[e]-1,3-oxazine, octahy...	216	C13H16N2O	1000138-92-2	14
2			2-Cyclohexen-3-ol-1-one, 2-benzoyl-	216	C13H12O3	061834-43-3	14
3			4-Methyl-2H,3H,3aH-pyrrolo[1,2-a...	216	C12H12N2O2	1000387-53-2	14
4			Succinic acid, 2,2,3,3-tetraflu...	370	C17H26F4O4	1000390-19-4	12
5			Silane, methylvinyl(1-cyclopenty...	328	C16H32O3Si2	1000420-98-4	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ281

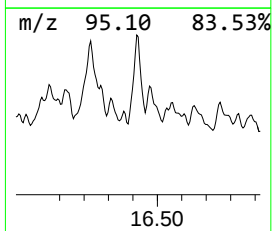
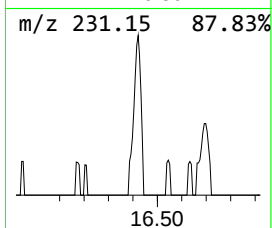
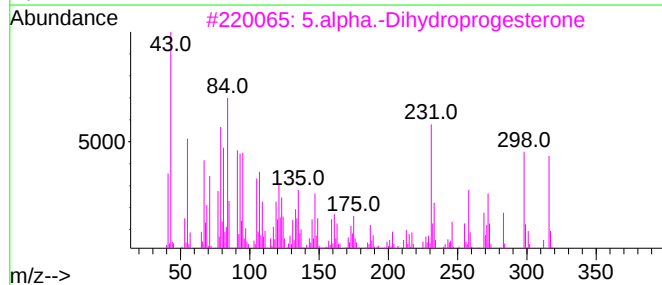
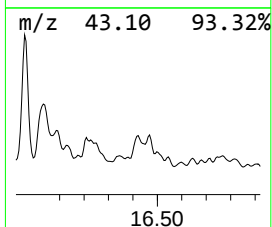
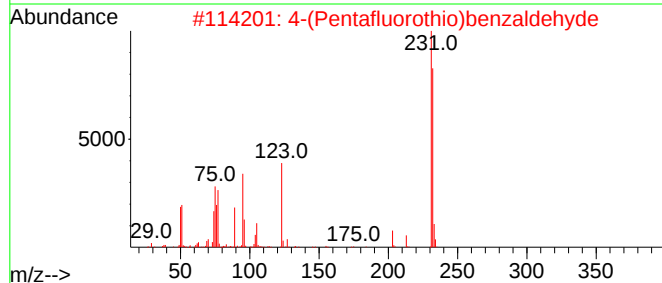
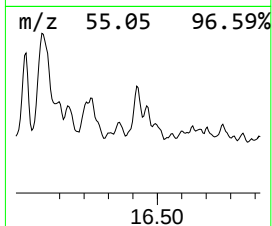
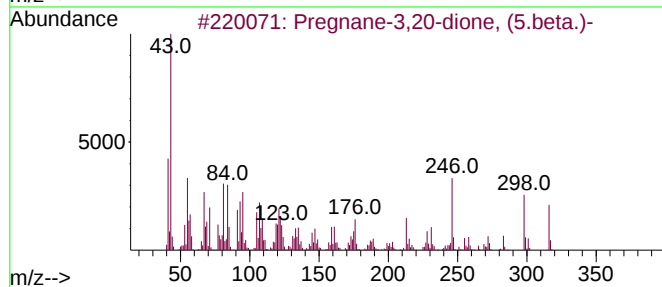
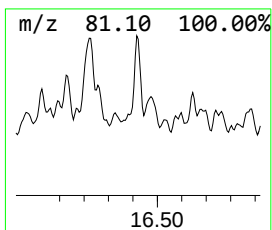
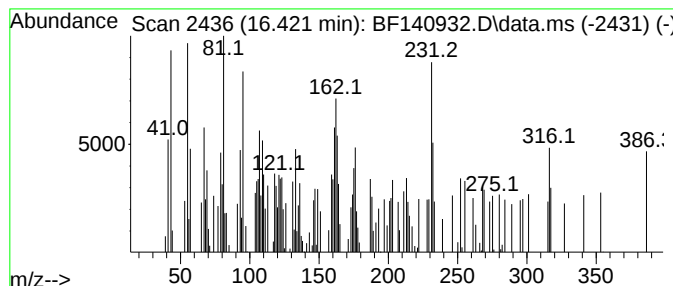
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.421 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.421	2.06 ng	38941	Perylene-d12	15.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pregnane-3,20-dione, (5.beta.)-	316	C21H32O2	000128-23-4	42
2			4-(Pentafluorothio)benzaldehyde	232	C7H5F5OS	401892-84-0	22
3			5.alpha.-Dihydroprogesterone	316	C21H32O2	000566-65-4	18
4			(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	15
5			Phosphinic acid, (2-methylphenyl)...	232	C13H13O2P	018593-18-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121924\
Data File : BF140932.D
Acq On : 19 Dec 2024 17:49
Operator : RC/JU
Sample : P5343-03
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ281

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.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
Benzophenone	10.592	5.0	ng	108139	3	9.881	428891	20.0
n-Hexadecanoic ...	11.892	10.5	ng	181419	4	11.374	346623	20.0
5-Bromo-2-thiop...	11.986	4.5	ng	77467	4	11.374	346623	20.0
unknown12.680	12.680	2.2	ng	38294	4	11.374	346623	20.0
Benzo[b]naphtho...	13.780	2.4	ng	84269	5	14.033	695474	20.0
Heptadecyl trif...	13.851	3.5	ng	120830	5	14.033	695474	20.0
1-Tetracosene	14.510	5.5	ng	192223	5	14.033	695474	20.0
1-Nonadecene	15.180	15.3	ng	289032	6	15.545	377818	20.0
Benzo[e]pyrene	15.416	5.0	ng	93919	6	15.545	377818	20.0
Octacosane, 2-m...	15.957	2.6	ng	48974	6	15.545	377818	20.0
Cyclopentadecane	16.033	4.6	ng	87684	6	15.545	377818	20.0
unknown16.221	16.221	3.8	ng	71085	6	15.545	377818	20.0
unknown16.421	16.421	2.1	ng	38941	6	15.545	377818	20.0