

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
 Data File : BF139578.D
 Acq On : 30 Sep 2024 15:00
 Operator : RC/JU
 Sample : P4213-01 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-02-092624

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-BF091324.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139578.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.057	495	504	512	rVB	21933	31807	3.68%	0.262%
2	5.475	569	575	583	rBV	476653	606885	70.17%	4.999%
3	6.487	742	747	756	rBV	499982	643140	74.36%	5.298%
4	6.863	806	811	816	rBV	330638	396632	45.86%	3.267%
5	7.422	901	906	911	rBV	337823	410184	47.43%	3.379%
6	7.681	946	950	957	rVB	10069	12889	1.49%	0.106%
7	8.145	1024	1029	1034	rBV	417263	498214	57.61%	4.104%
8	9.216	1206	1211	1216	rBV	588997	699596	80.89%	5.763%
9	9.292	1220	1224	1228	rBV	9266	12182	1.41%	0.100%
10	9.557	1262	1269	1271	rBV4	6421	10314	1.19%	0.085%
11	9.610	1275	1278	1282	rBV3	5910	8736	1.01%	0.072%
12	9.822	1307	1314	1318	rVV	14973	22746	2.63%	0.187%
13	9.898	1321	1327	1330	rVV	422496	538362	62.25%	4.435%
14	9.928	1330	1332	1336	rVB	39353	41715	4.82%	0.344%
15	10.098	1356	1361	1365	rVV	20742	32327	3.74%	0.266%
16	10.139	1365	1368	1373	rVV6	7548	11720	1.36%	0.097%
17	10.322	1392	1399	1403	rBV	21390	29923	3.46%	0.247%
18	10.439	1415	1419	1425	rBV	29237	41129	4.76%	0.339%
19	10.510	1427	1431	1432	rVV3	9914	13148	1.52%	0.108%
20	10.539	1433	1436	1441	rVV	15146	27578	3.19%	0.227%
21	10.604	1442	1447	1454	rVV	120794	156424	18.09%	1.289%
22	10.681	1455	1460	1471	rVV	201236	262149	30.31%	2.160%
23	10.798	1475	1480	1488	rVB	29262	50421	5.83%	0.415%
24	10.916	1497	1500	1505	rVB	7969	9468	1.09%	0.078%
25	10.986	1508	1512	1515	rBV4	8318	11665	1.35%	0.096%
26	11.122	1530	1535	1537	rBV5	5972	10340	1.20%	0.085%
27	11.186	1542	1546	1550	rVV	11274	14798	1.71%	0.122%
28	11.239	1550	1555	1558	rVV2	23614	35323	4.08%	0.291%
29	11.275	1558	1561	1567	rVB2	24973	33816	3.91%	0.279%
30	11.380	1574	1579	1581	rBV	381586	501926	58.04%	4.135%
31	11.404	1581	1583	1587	rVV	238294	257003	29.72%	2.117%
32	11.451	1587	1591	1595	rVV	56071	75942	8.78%	0.626%
33	11.610	1611	1618	1624	rBV2	26953	50210	5.81%	0.414%
34	11.669	1624	1628	1632	rVV2	21998	31950	3.69%	0.263%
35	11.710	1632	1635	1641	rVB4	9271	14041	1.62%	0.116%
36	11.904	1660	1668	1674	rBV2	214326	374849	43.34%	3.088%

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37	11.992	1679	1683	1691	rVV	70453	126796	14.66%	1.045%
38	12.075	1693	1697	1701	rVV2	18580	26580	3.07%	0.219%
39	12.175	1711	1714	1716	rBV	17850	22393	2.59%	0.184%
40	12.198	1716	1718	1725	rVB3	23425	29502	3.41%	0.243%
41	12.351	1742	1744	1747	rBV	9835	12966	1.50%	0.107%
42	12.427	1753	1757	1760	rBV	21532	30995	3.58%	0.255%
43	12.463	1760	1763	1768	rVV2	32175	47189	5.46%	0.389%
44	12.516	1768	1772	1775	rVB	18119	24283	2.81%	0.200%
45	12.592	1780	1785	1790	rBV	503595	632804	73.17%	5.213%
46	12.692	1797	1802	1817	rBV	471772	864852	100.00%	7.125%
47	12.822	1817	1824	1829	rVB	428088	655062	75.74%	5.396%
48	12.957	1841	1847	1855	rVB	387844	564913	65.32%	4.654%
49	13.033	1856	1860	1864	rBV2	28121	50549	5.84%	0.416%
50	13.145	1872	1879	1883	rVB	61683	108601	12.56%	0.895%
51	13.192	1883	1887	1888	rBV3	22856	30748	3.56%	0.253%
52	13.210	1888	1890	1893	rVB	29402	30525	3.53%	0.251%
53	13.251	1893	1897	1899	rBV2	33352	43078	4.98%	0.355%
54	13.339	1910	1912	1915	rBV	12704	13436	1.55%	0.111%
55	13.374	1915	1918	1921	rBV	16302	18468	2.14%	0.152%
56	13.533	1941	1945	1948	rBV3	28755	46246	5.35%	0.381%
57	13.651	1963	1965	1970	rVB	29221	35800	4.14%	0.295%
58	13.763	1980	1984	1987	rBV3	41159	64062	7.41%	0.528%
59	13.857	1997	2000	2009	rVB	66636	106777	12.35%	0.880%
60	14.010	2020	2026	2030	rBV2	458925	812470	93.94%	6.693%
61	14.121	2042	2045	2048	rVB	39572	44512	5.15%	0.367%
62	14.480	2103	2106	2110	rBV3	38630	55255	6.39%	0.455%
63	14.786	2155	2158	2161	rBV	38822	46513	5.38%	0.383%
64	15.051	2198	2203	2211	rBV	193068	425881	49.24%	3.508%
65	15.351	2251	2254	2260	rVB	102757	161748	18.70%	1.332%
66	15.415	2261	2265	2271	rVV	142638	228512	26.42%	1.882%
67	15.480	2271	2276	2291	rVB2	196464	432437	50.00%	3.562%
68	16.180	2391	2395	2409	rVB2	73051	151541	17.52%	1.248%
69	16.933	2519	2523	2532	rVB2	56195	119804	13.85%	0.987%
70	17.374	2594	2598	2607	rVB	45344	94154	10.89%	0.776%

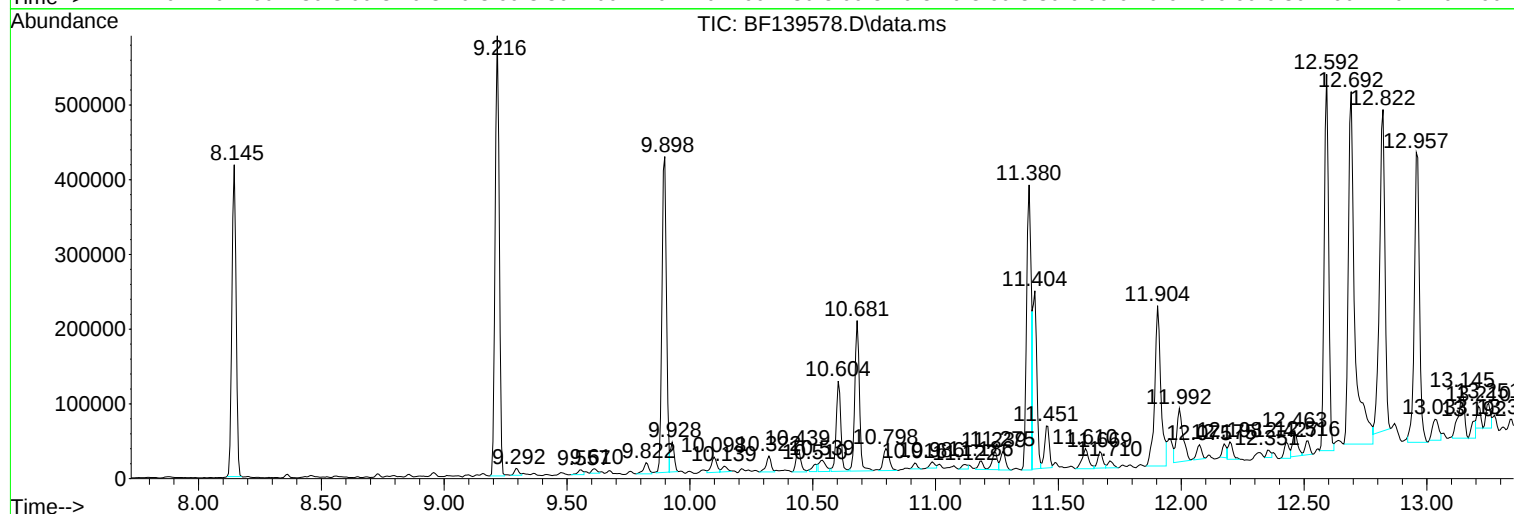
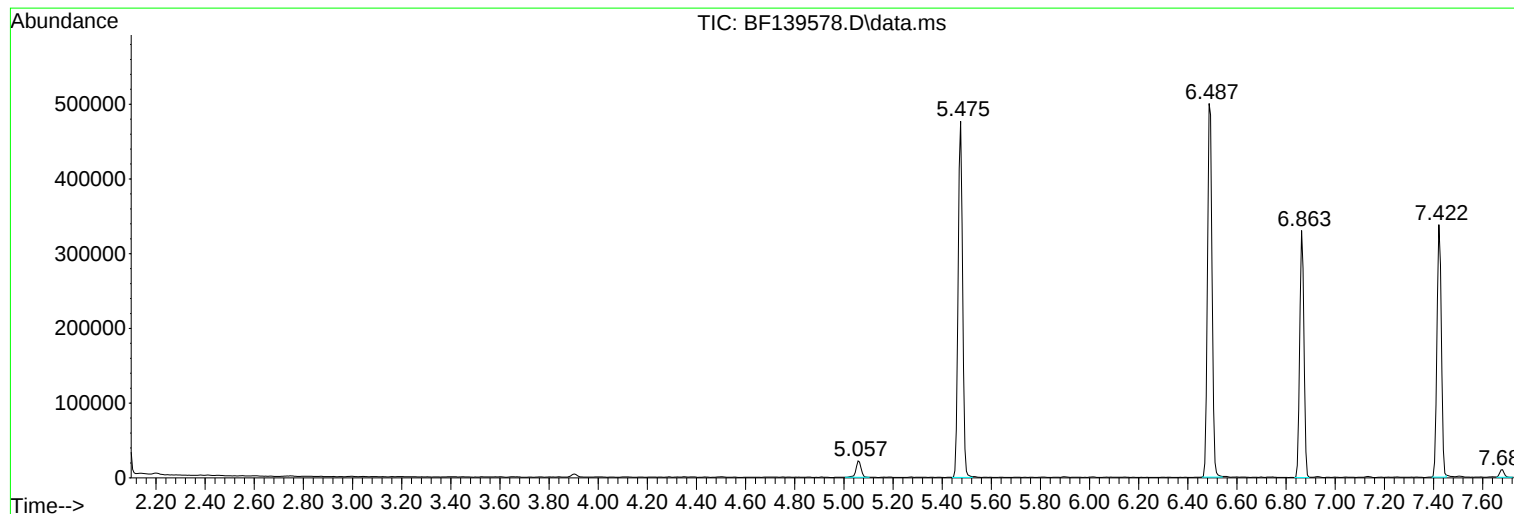
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : P4213-01 2X
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Instrument :
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HD-02-092624

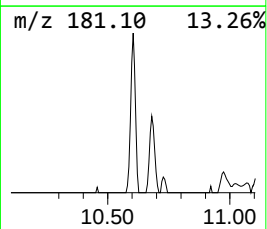
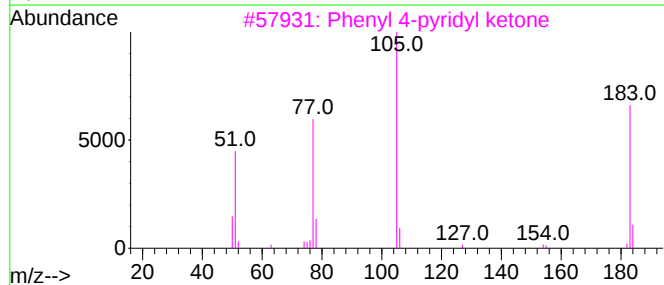
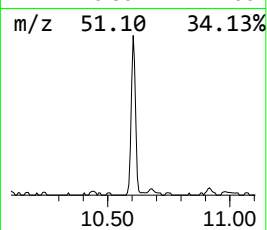
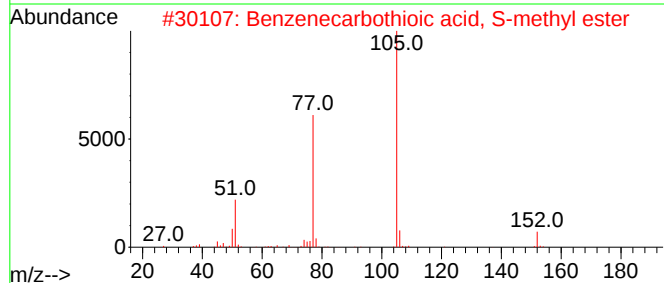
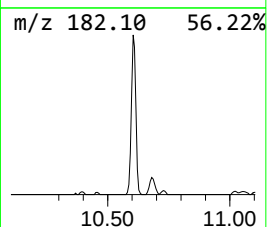
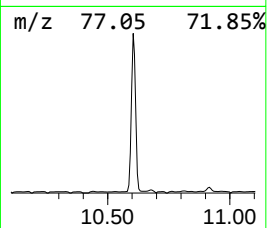
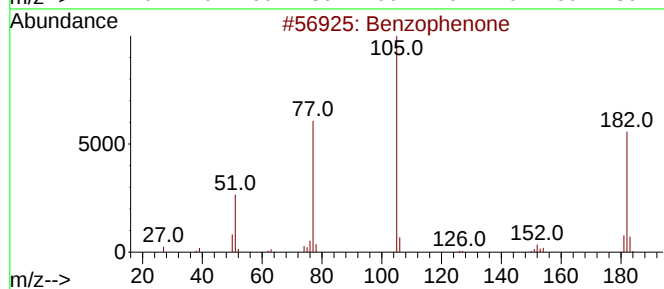
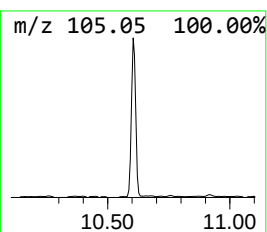
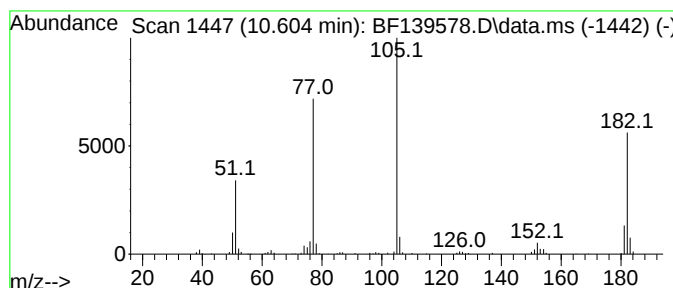
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.604	5.81 ng	156424	Acenaphthene-d10	9.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	60
3			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	53
4			N-cyanobenzamide	146	C8H6N2O	015150-25-1	49
5			2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	47



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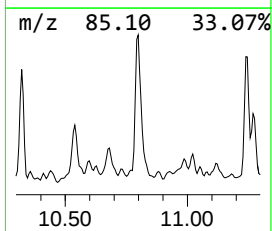
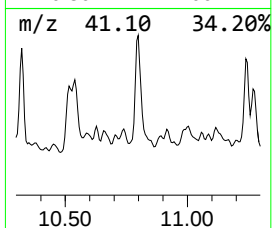
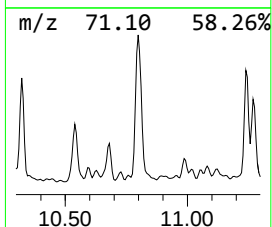
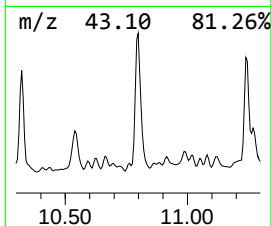
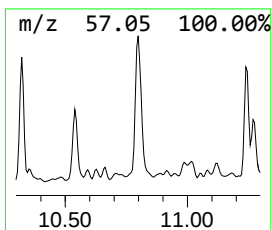
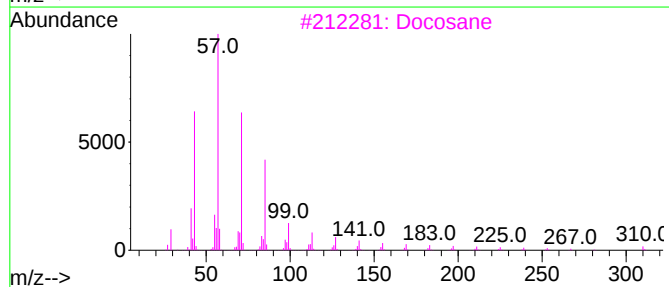
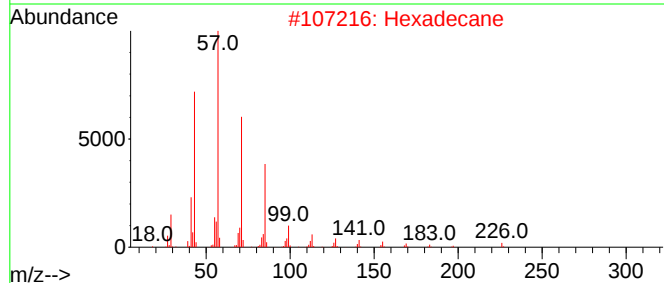
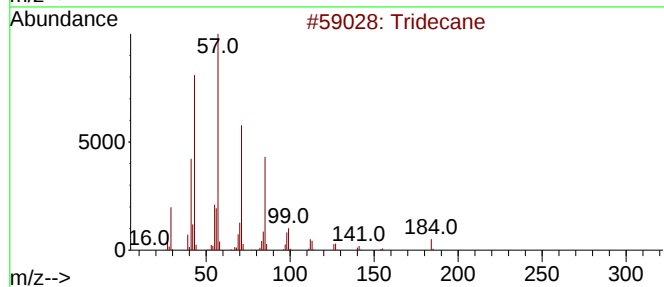
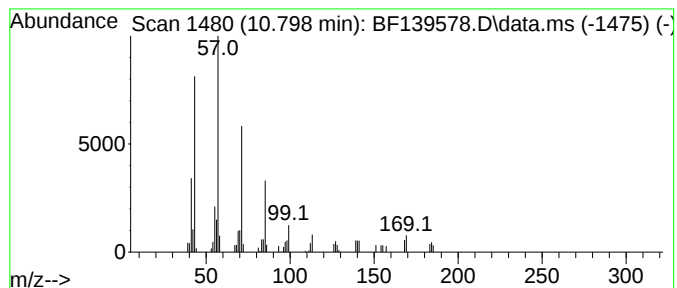
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Tridecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.798	2.01 ng	50421	Phenanthrene-d10	11.381

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	90
2		Hexadecane	226	C16H34	000544-76-3	80
3		Docosane	310	C22H46	000629-97-0	80
4		Eicosane	282	C20H42	000112-95-8	80
5		Octadecane	254	C18H38	000593-45-3	74



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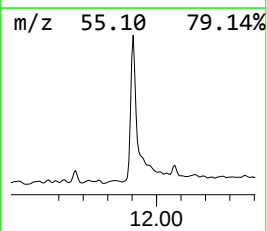
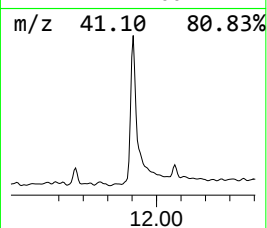
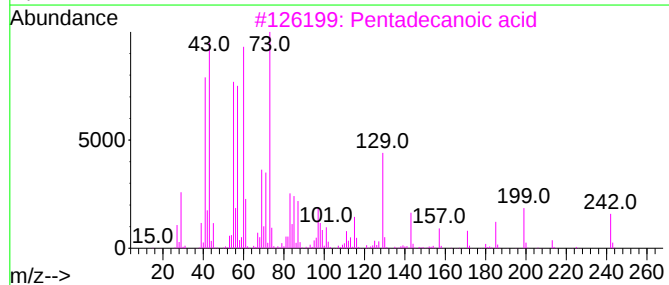
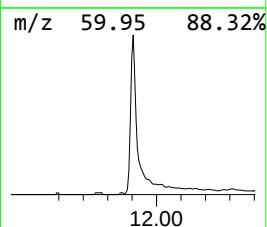
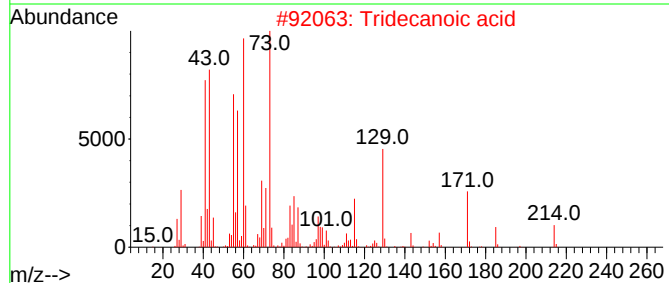
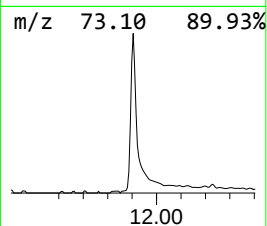
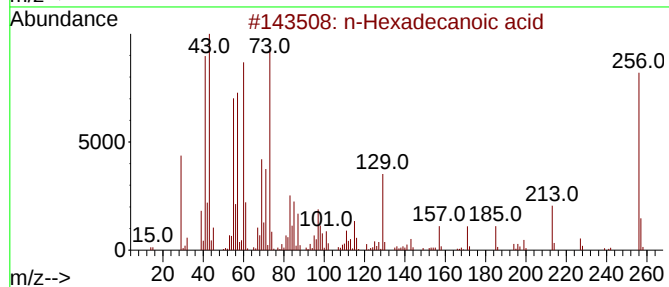
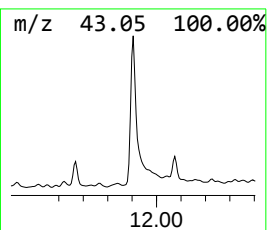
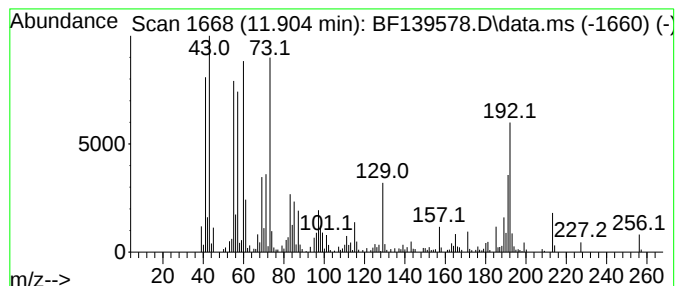
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.904	14.94 ng	374849	Phenanthrene-d10	11.381

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	86
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	76
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	70
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	59



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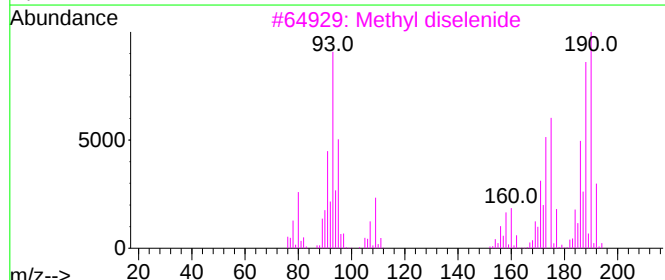
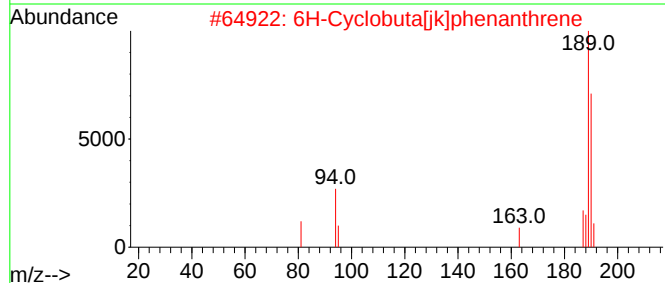
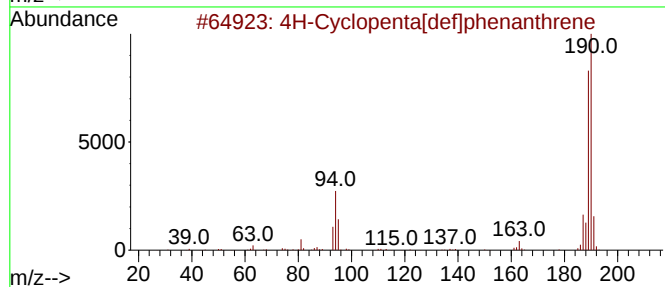
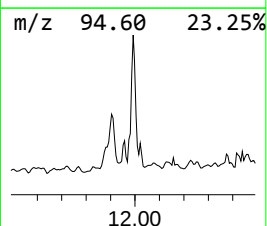
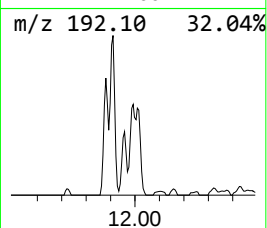
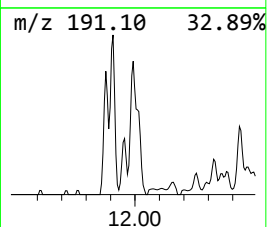
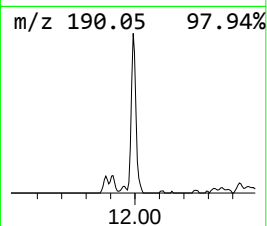
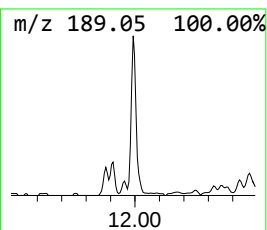
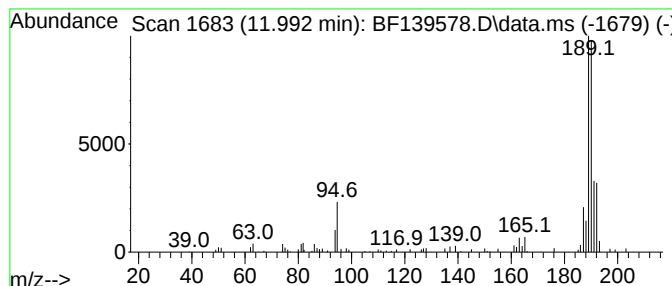
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.992	5.05 ng	126796	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3			Methyl diselenide	190	C2H6Se2	007101-31-7	46
4			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
5			1,2,3,4,5,6,7,8,9,10-Decahydro-5...	190	C13H18O	1000497-99-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
Data File : BF139578.D
Acq On : 30 Sep 2024 15:00
Operator : RC/JU
Sample : P4213-01 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-092624

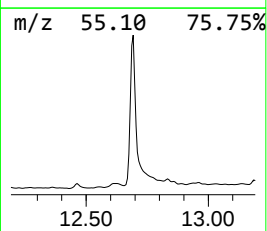
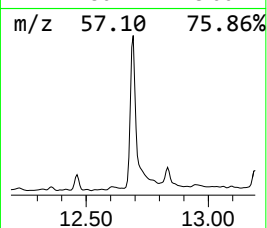
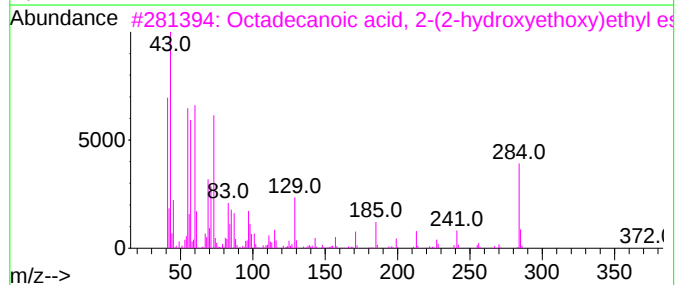
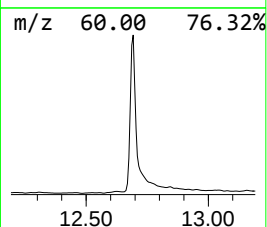
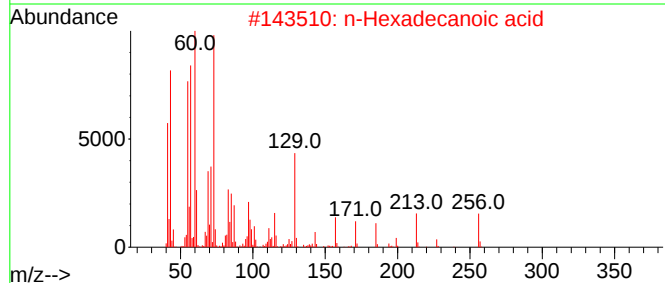
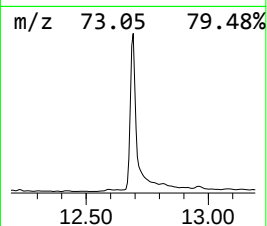
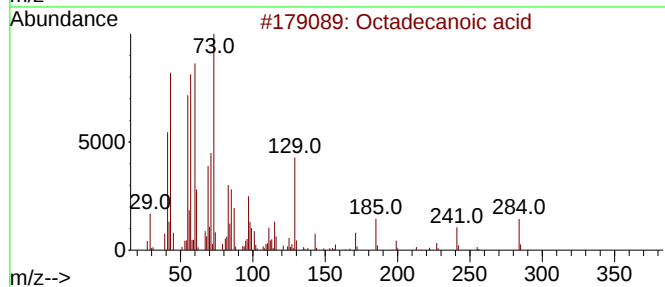
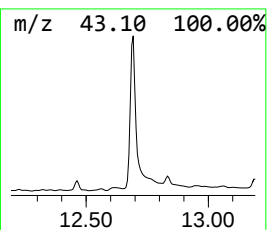
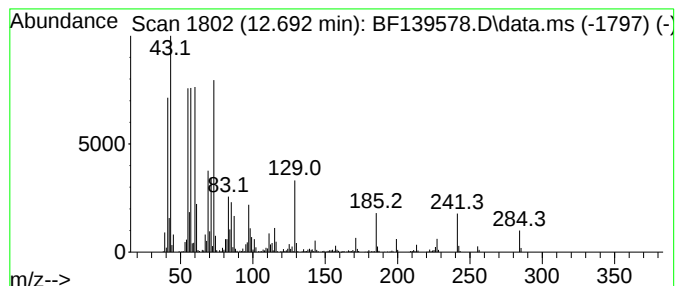
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.692	34.46 ng	864852	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
4			Isopropyl palmitate	298	C19H38O2	000142-91-6	72
5			Tridecanoic acid	214	C13H26O2	000638-53-9	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
Data File : BF139578.D
Acq On : 30 Sep 2024 15:00
Operator : RC/JU
Sample : P4213-01 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-092624

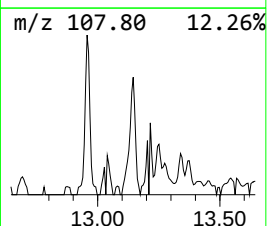
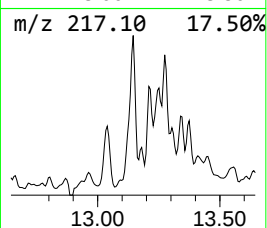
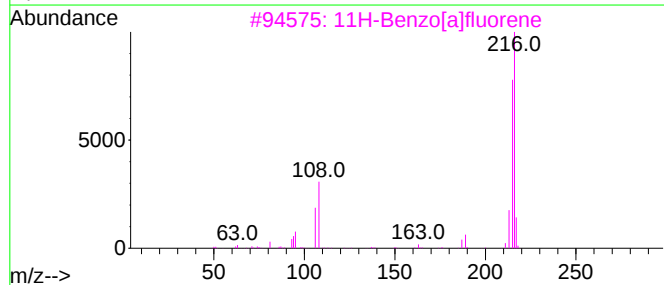
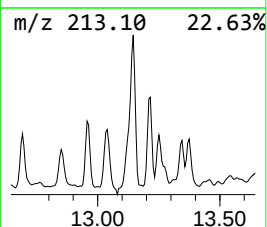
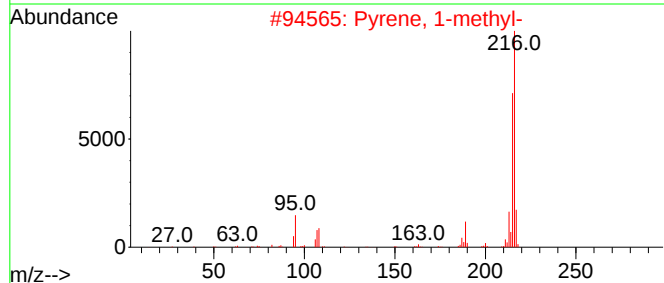
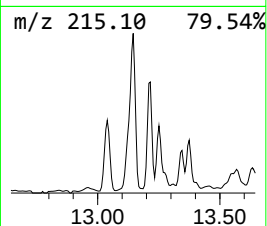
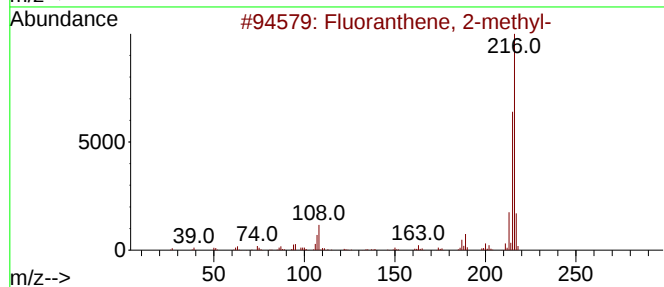
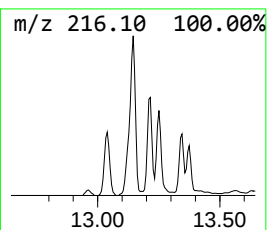
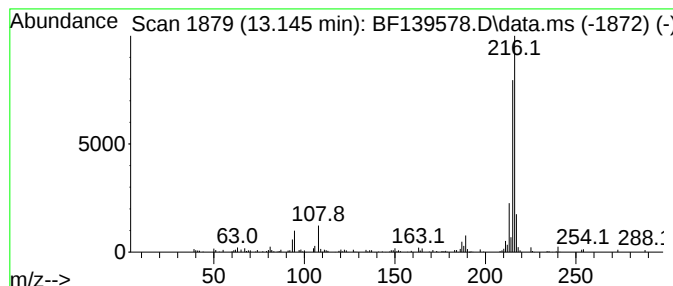
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.145	2.67 ng	108601	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
Data File : BF139578.D
Acq On : 30 Sep 2024 15:00
Operator : RC/JU
Sample : P4213-01 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-092624

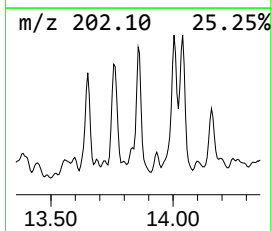
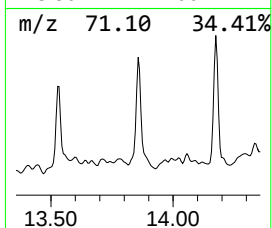
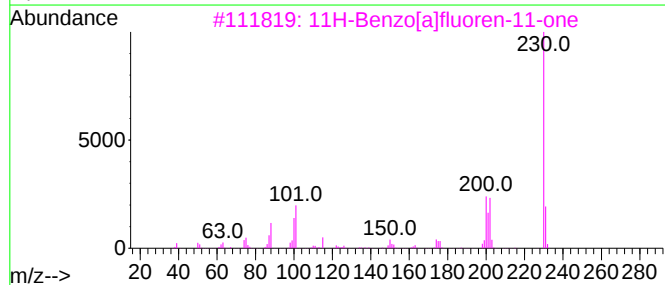
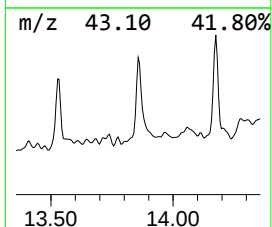
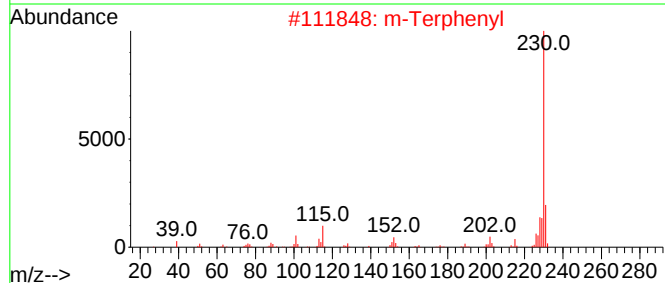
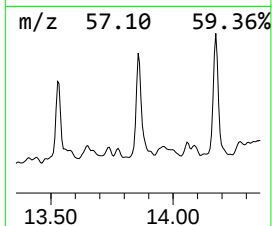
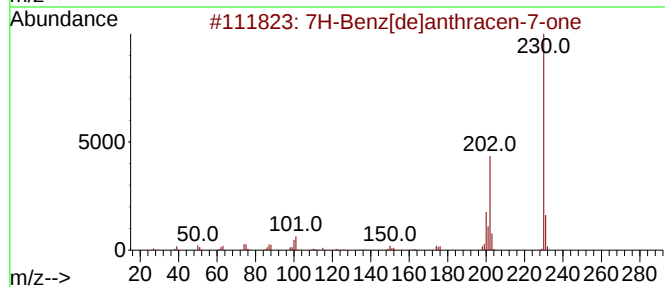
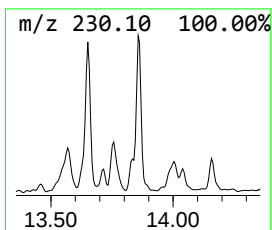
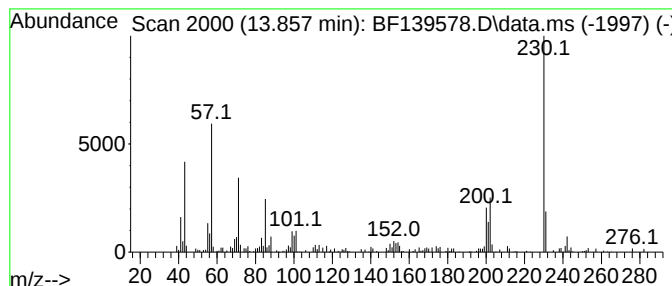
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 7H-Benz[de]anthracen-7-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	2.63 ng	106777	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
2			m-Terphenyl	230	C18H14	000092-06-8	43
3			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	38
4			1,1'-Biphenyl, 5-hydroxy-3,4'-di...	230	C14H14O3	119101-26-7	35
5			p-Terphenyl	230	C18H14	000092-94-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
Data File : BF139578.D
Acq On : 30 Sep 2024 15:00
Operator : RC/JU
Sample : P4213-01 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-092624

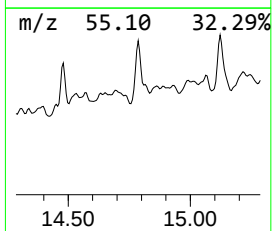
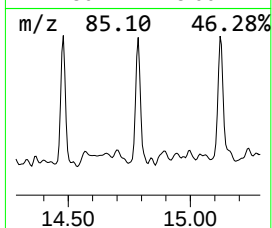
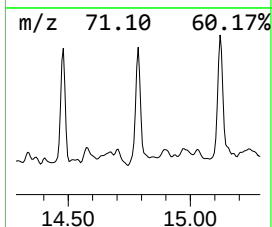
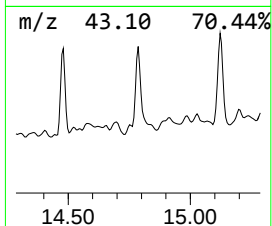
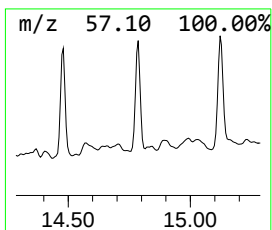
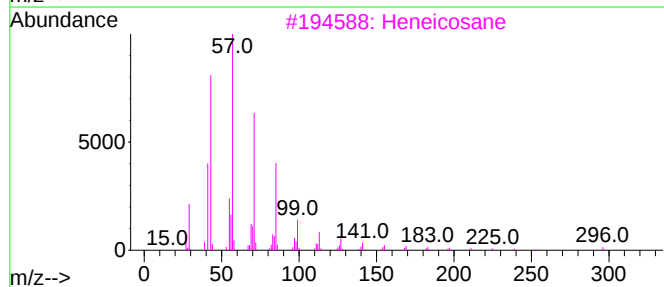
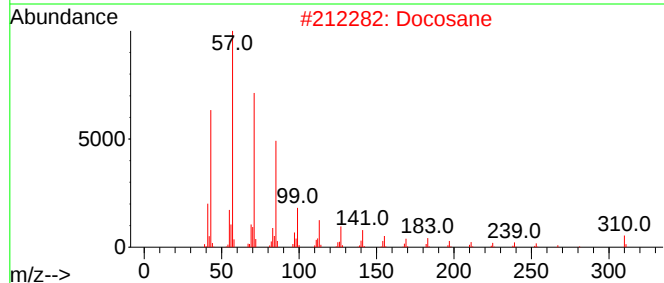
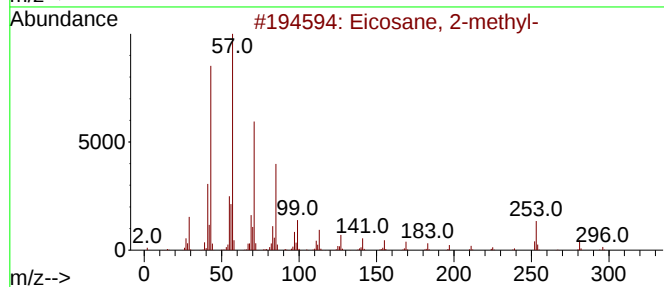
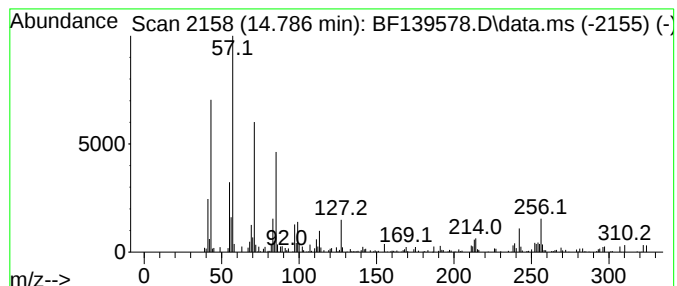
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.786	2.15 ng	46513	Perylene-d12	15.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 2-methyl-	296	C21H44	001560-84-5	90
2			Docosane	310	C22H46	000629-97-0	86
3			Heneicosane	296	C21H44	000629-94-7	86
4			Pentadecane	212	C15H32	000629-62-9	86
5			Octadecane	254	C18H38	000593-45-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
Data File : BF139578.D
Acq On : 30 Sep 2024 15:00
Operator : RC/JU
Sample : P4213-01 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-092624

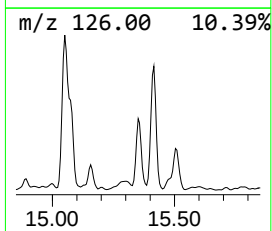
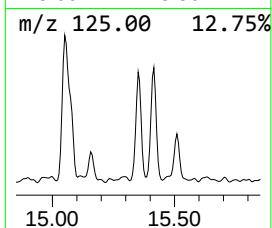
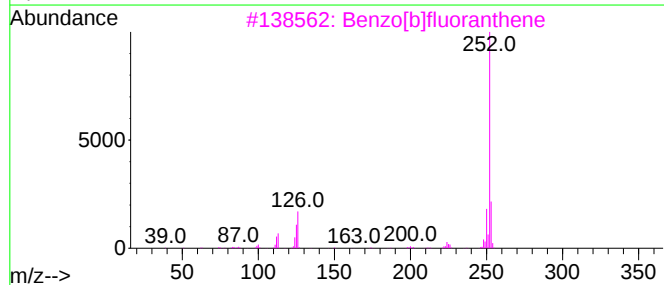
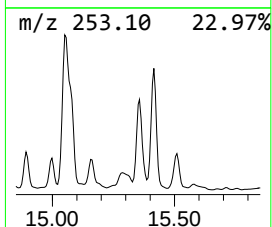
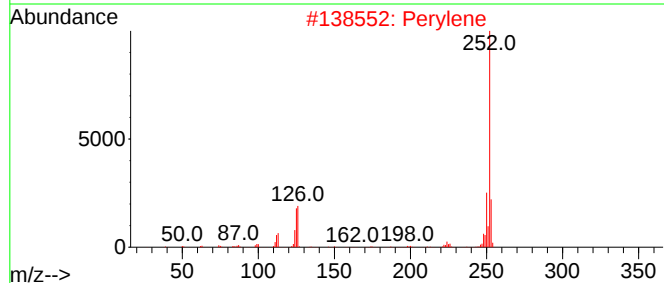
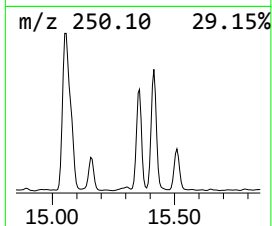
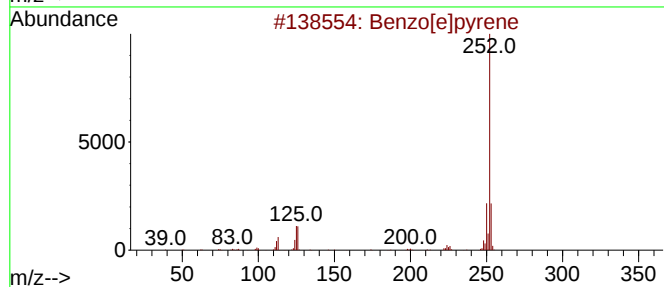
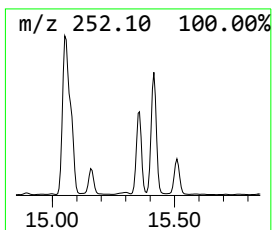
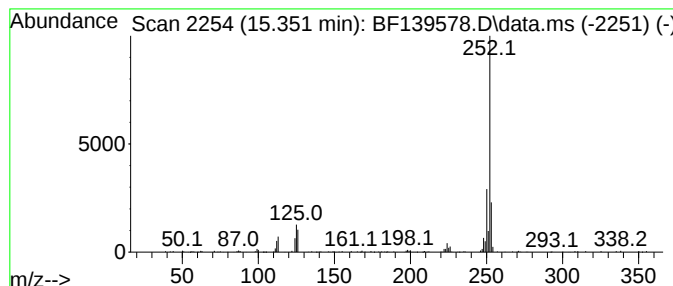
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.351	7.48 ng	161748	Perylene-d12	15.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Perylene	252	C20H12	000198-55-0	94
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
Data File : BF139578.D
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Misc :
ALS Vial : 10 Sample Multiplier: 1

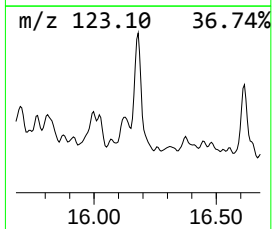
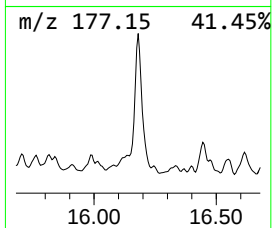
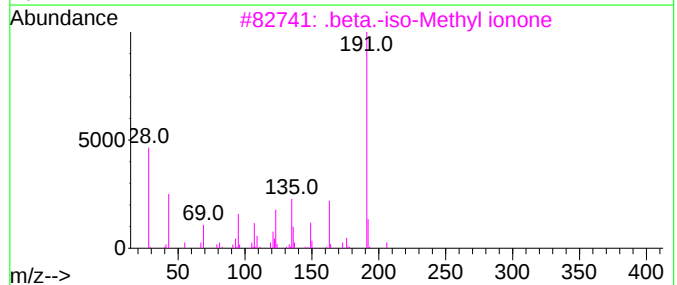
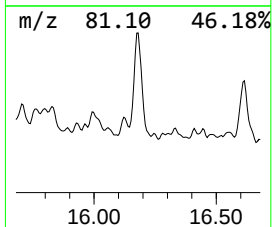
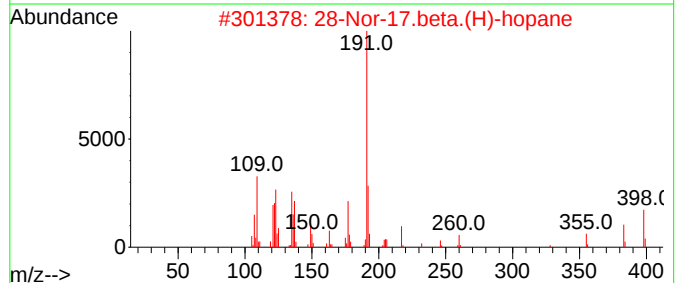
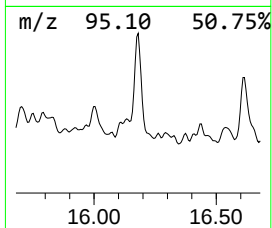
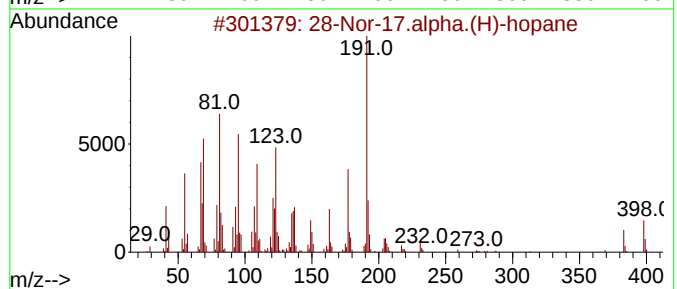
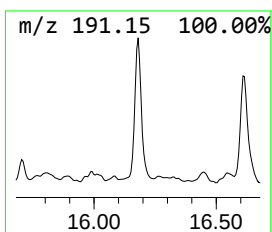
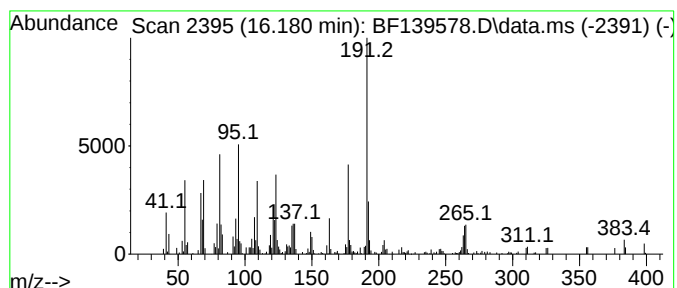
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 28-Nor-17.alpha.(H)-hopane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.180	7.01 ng	151541	Perylene-d12	15.480	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	70
2	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	64
3	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	45
4	Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	35
5	(3-Cyano-6-ethyl-5-methyl-pyridi...	264	C13H16N2O2S	1000300-45-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093024\
Data File : BF139578.D
Acq On : 30 Sep 2024 15:00
Operator : RC/JU
Sample : P4213-01 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HD-02-092624

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.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
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Tridecane	10.798	2.0	ng	50421	4	11.381	501926	20.0
n-Hexadecanoic ...	11.904	14.9	ng	374849	4	11.381	501926	20.0
4H-Cyclopenta[d...]	11.992	5.0	ng	126796	4	11.381	501926	20.0
Octadecanoic acid	12.692	34.5	ng	864852	4	11.381	501926	20.0
Fluoranthene, 2...	13.145	2.7	ng	108601	5	14.016	812470	20.0
7H-Benz[de]anth...	13.857	2.6	ng	106777	5	14.016	812470	20.0
Eicosane, 2-met...	14.786	2.1	ng	46513	6	15.480	432437	20.0
Benzo[e]pyrene	15.351	7.5	ng	161748	6	15.480	432437	20.0
28-Nor-17.alpha...	16.180	7.0	ng	151541	6	15.480	432437	20.0