

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
 Data File : BF137479.D
 Acq On : 04 Mar 2024 14:20
 Operator : CG\JU
 Sample : P1662-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-352

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-BF022924.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF137479.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.187	11	17	33	rVB	1380340	1918679	38.63%	3.103%
2	5.134	514	518	524	rBV	42887	59105	1.19%	0.096%
3	5.516	579	583	609	rBV	4420925	4401510	88.62%	7.118%
4	6.498	746	750	753	rBV	4223904	4369642	87.98%	7.066%
5	6.857	808	811	824	rBV	1440442	1310499	26.39%	2.119%
6	7.416	902	906	918	rBV	2780917	2853135	57.45%	4.614%
7	8.133	1024	1028	1048	rBV	1875857	1786196	35.97%	2.888%
8	8.457	1077	1083	1088	rBV	32614	50146	1.01%	0.081%
9	8.851	1146	1150	1156	rBV	58225	62277	1.25%	0.101%
10	8.945	1162	1166	1174	rVB	115988	110410	2.22%	0.179%
11	9.210	1206	1211	1215	rBV	4979877	4957880	99.83%	8.017%
12	9.316	1222	1229	1235	rVB	121550	144490	2.91%	0.234%
13	9.474	1252	1256	1265	rBV	119323	175982	3.54%	0.285%
14	9.551	1265	1269	1271	rBV	176460	184904	3.72%	0.299%
15	9.574	1271	1273	1279	rVB	134332	113090	2.28%	0.183%
16	9.669	1286	1289	1295	rBV	73179	96276	1.94%	0.156%
17	9.751	1295	1303	1308	rVB	57543	60751	1.22%	0.098%
18	9.892	1320	1327	1329	rBV	2416205	2136568	43.02%	3.455%
19	9.922	1329	1332	1338	rVV	1646934	1405058	28.29%	2.272%
20	9.998	1341	1345	1350	rVV	97901	148719	2.99%	0.240%
21	10.051	1350	1354	1356	rVV2	75054	77897	1.57%	0.126%
22	10.098	1358	1362	1366	rVV	927942	834828	16.81%	1.350%
23	10.133	1366	1368	1375	rVB	67795	83019	1.67%	0.134%
24	10.304	1392	1397	1398	rBV	51543	54200	1.09%	0.088%
25	10.439	1417	1420	1426	rBV	1454349	1248970	25.15%	2.020%
26	10.527	1433	1435	1438	rVB	159040	120351	2.42%	0.195%
27	10.598	1445	1447	1452	rVB	250806	244862	4.93%	0.396%
28	10.686	1458	1462	1468	rBV	3397612	3225925	64.95%	5.217%
29	10.733	1468	1470	1474	rVB	67810	64528	1.30%	0.104%
30	10.816	1478	1484	1488	rVB3	92031	115076	2.32%	0.186%
31	10.992	1508	1514	1517	rBV2	175479	239576	4.82%	0.387%
32	11.074	1522	1528	1531	rVB2	71491	91939	1.85%	0.149%
33	11.121	1531	1536	1538	rBV2	72829	79523	1.60%	0.129%
34	11.192	1544	1548	1552	rVB2	172018	178138	3.59%	0.288%
35	11.233	1552	1555	1561	rVV3	83708	142248	2.86%	0.230%
36	11.280	1561	1563	1571	rVB	296829	295021	5.94%	0.477%

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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-BF022924.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.386	1577	1581	1583	rBV	2268300	2151307	43.32%	3.479%
38	11.416	1583	1586	1589	rVV	4533380	4151828	83.60%	6.714%
39	11.463	1589	1594	1597	rVB	393095	410007	8.26%	0.663%
40	11.621	1618	1621	1630	rBV2	125516	221570	4.46%	0.358%
41	11.686	1630	1632	1636	rBV	99971	85005	1.71%	0.137%
42	11.804	1646	1652	1657	rVB3	37216	56865	1.14%	0.092%
43	11.892	1664	1667	1669	rBV	371239	303900	6.12%	0.491%
44	11.921	1669	1672	1673	rVV	515508	528046	10.63%	0.854%
45	11.933	1673	1674	1678	rVV	646481	594019	11.96%	0.961%
46	11.968	1678	1680	1683	rVV	210342	229757	4.63%	0.372%
47	12.004	1683	1686	1689	rVV	901586	875033	17.62%	1.415%
48	12.027	1689	1690	1695	rVB	174555	109732	2.21%	0.177%
49	12.192	1715	1718	1721	rBV	231649	207401	4.18%	0.335%
50	12.215	1721	1722	1727	rVB	66741	51223	1.03%	0.083%
51	12.339	1735	1743	1746	rBV3	72050	116232	2.34%	0.188%
52	12.445	1757	1761	1764	rBV	86015	96293	1.94%	0.156%
53	12.486	1764	1768	1773	rVB4	74125	116388	2.34%	0.188%
54	12.533	1773	1776	1780	rBV2	67386	78035	1.57%	0.126%
55	12.610	1785	1789	1792	rBV	3344933	3000968	60.42%	4.853%
56	12.721	1805	1808	1812	rBV2	104331	135493	2.73%	0.219%
57	12.757	1812	1814	1822	rVV	91133	103313	2.08%	0.167%
58	12.839	1822	1828	1831	rVV2	2161937	2518695	50.71%	4.073%
59	12.892	1834	1837	1841	rVB2	105407	123926	2.50%	0.200%
60	12.963	1845	1849	1850	rBV	111965	109803	2.21%	0.178%
61	12.992	1850	1854	1857	rVV	5204874	4966461	100.00%	8.031%
62	13.063	1860	1866	1870	rVB2	90854	156564	3.15%	0.253%
63	13.145	1877	1880	1882	rBV2	78254	91665	1.85%	0.148%
64	13.168	1882	1884	1888	rVB2	282029	247053	4.97%	0.400%
65	13.233	1891	1895	1899	rBV2	284225	307214	6.19%	0.497%
66	13.274	1899	1902	1904	rBV	115360	107211	2.16%	0.173%
67	13.398	1921	1923	1932	rVB4	49664	60000	1.21%	0.097%
68	13.498	1936	1940	1944	rVB3	39243	58788	1.18%	0.095%
69	13.574	1948	1953	1955	rBV4	82643	81351	1.64%	0.132%
70	13.792	1987	1990	1993	rBV	69720	76987	1.55%	0.124%
71	13.821	1993	1995	1997	rBV	72043	65930	1.33%	0.107%
72	13.904	2005	2009	2012	rBV2	237346	289572	5.83%	0.468%
73	13.933	2012	2014	2018	rVB	85286	86513	1.74%	0.140%
74	14.045	2026	2033	2036	rBV2	1780865	2063140	41.54%	3.336%
75	14.074	2036	2038	2046	rVV	369543	412495	8.31%	0.667%
76	14.151	2048	2051	2054	rVB	88421	85057	1.71%	0.138%

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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-BF022924.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	14.227	2061	2064	2067	rVB	81254	71802	1.45%	0.116%
78	14.533	2113	2116	2119	rVB2	93991	86656	1.74%	0.140%
79	14.862	2168	2172	2177	rVB	102810	130357	2.62%	0.211%
80	14.939	2182	2185	2189	rBV	106008	130789	2.63%	0.212%
81	15.121	2211	2216	2219	rBV	133294	211643	4.26%	0.342%
82	15.233	2229	2235	2240	rVB2	173047	231158	4.65%	0.374%
83	15.439	2266	2270	2277	rVV	56719	82510	1.66%	0.133%
84	15.504	2277	2281	2288	rVV2	80663	144098	2.90%	0.233%
85	15.574	2288	2293	2301	rVV	885729	1247050	25.11%	2.017%
86	15.645	2301	2305	2313	rVB	102192	170710	3.44%	0.276%
87	16.133	2380	2388	2394	rBV	97534	202221	4.07%	0.327%
88	16.698	2479	2484	2493	rVB2	44390	93733	1.89%	0.152%
89	17.368	2593	2598	2607	rBV3	29318	63688	1.28%	0.103%

Sum of corrected areas: 61838673

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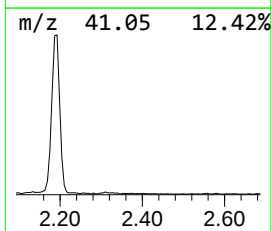
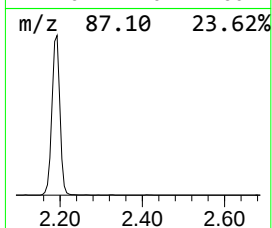
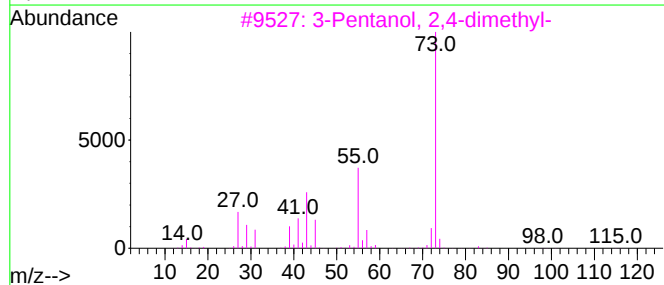
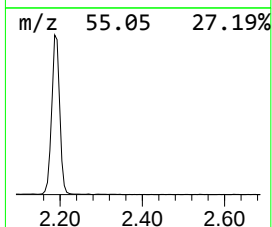
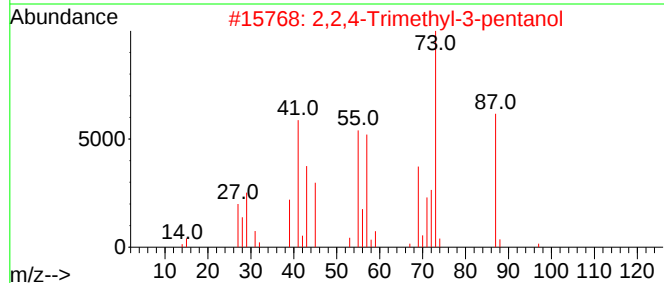
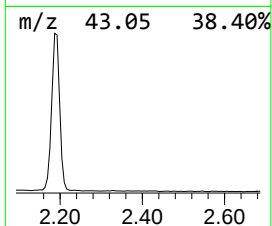
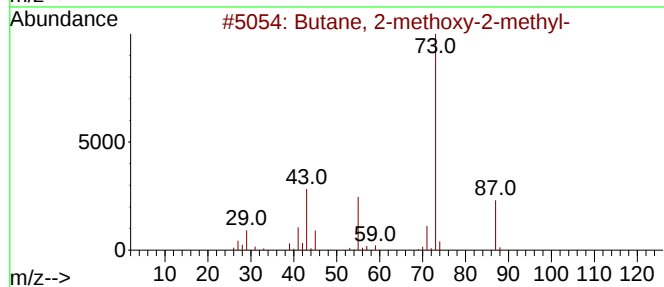
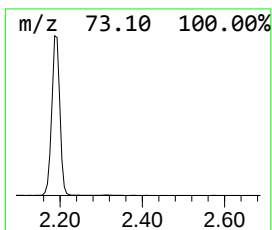
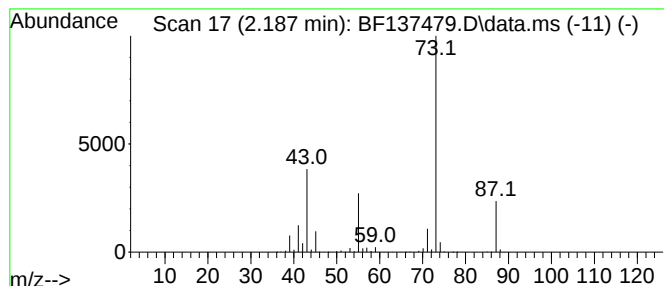
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.187	29.28 ng	1918680	1,4-Dichlorobenzene-d4	6.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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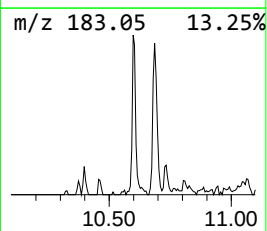
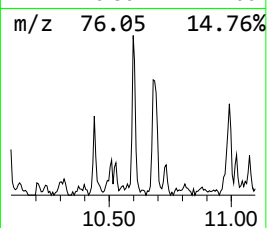
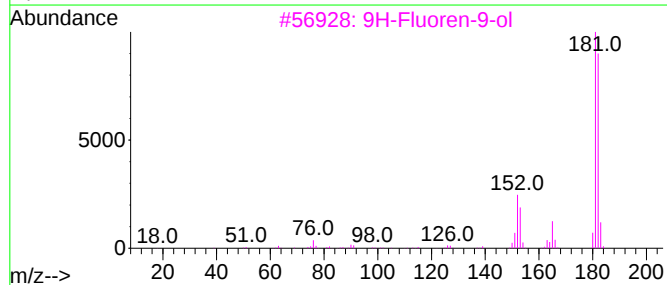
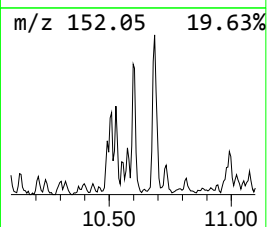
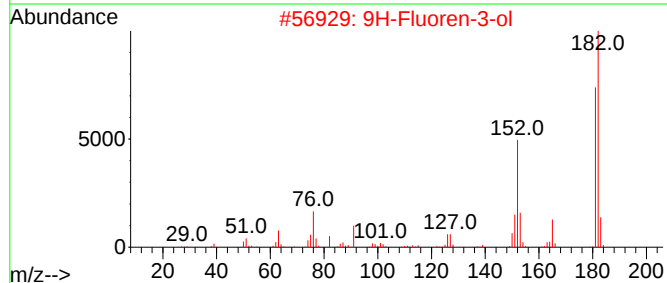
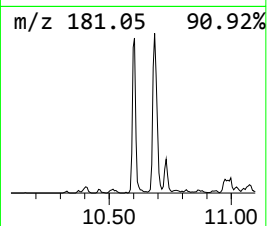
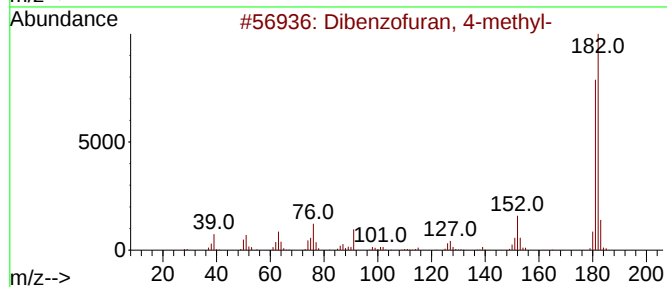
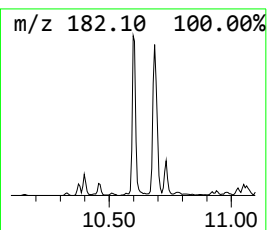
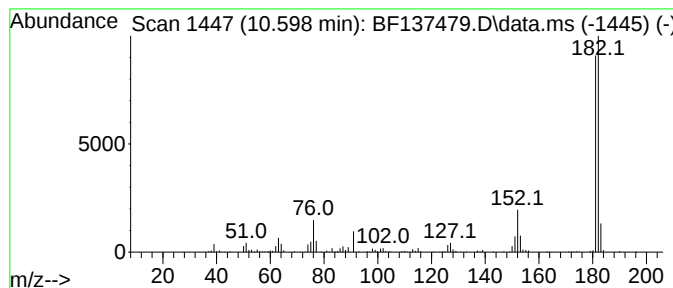
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Dibenzofuran, 4-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.598	2.29 ng	244862	Acenaphthene-d10	9.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	94
2			9H-Fluoren-3-ol	182	C13H10O	006344-67-8	91
3			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	72
4			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	72
5			2-Hydroxyfluorene	182	C13H10O	002443-58-5	64



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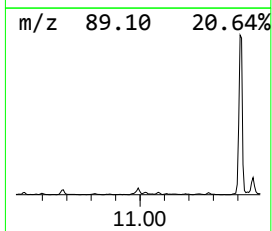
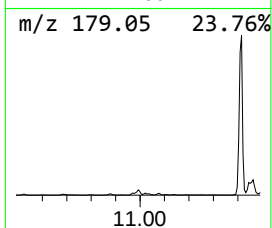
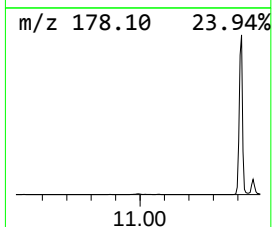
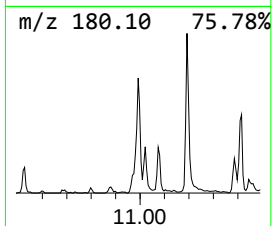
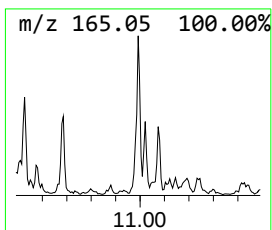
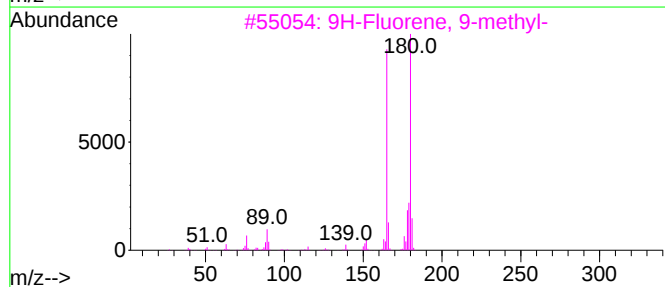
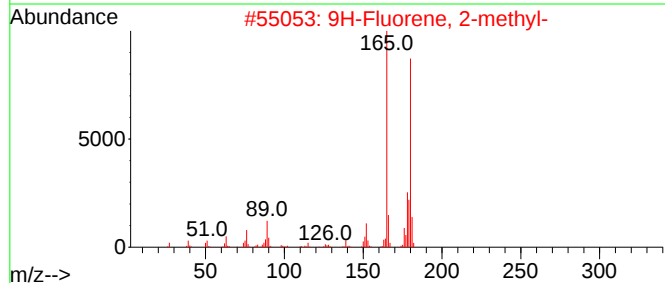
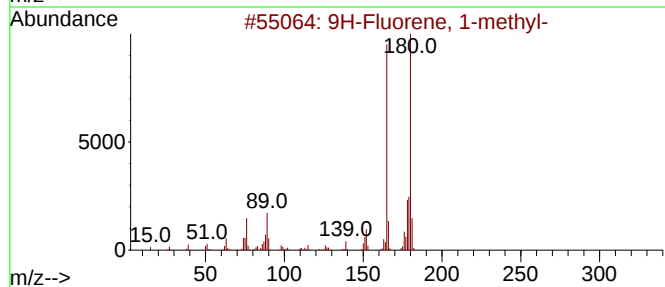
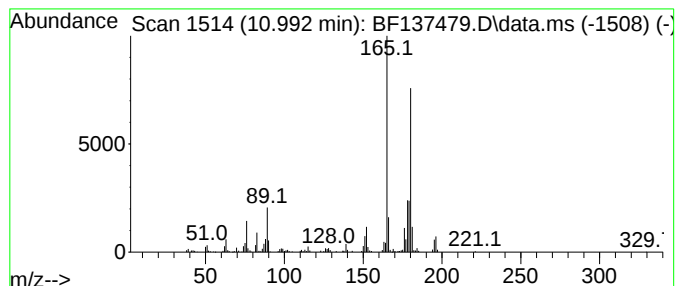
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 9H-Fluorene, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.992	2.23 ng	239576	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 1-methyl-			180	C14H12	001730-37-6	96
2	9H-Fluorene, 2-methyl-			180	C14H12	001430-97-3	96
3	9H-Fluorene, 9-methyl-			180	C14H12	002523-37-7	91
4	9H-Fluorene, 3-methyl-			180	C14H12	002523-39-9	90
5	9H-Fluorene, 4-methyl-			180	C14H12	001556-99-6	81



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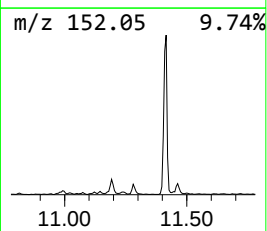
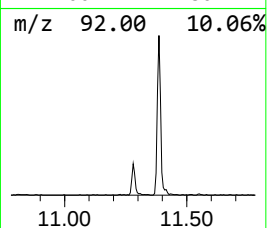
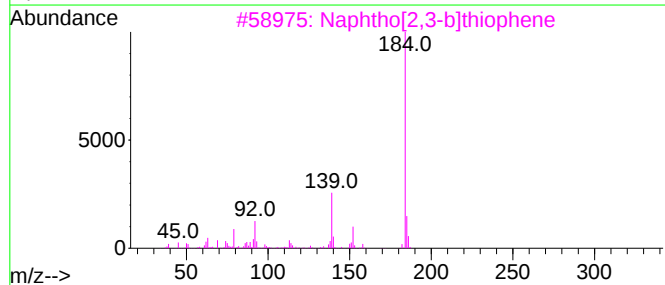
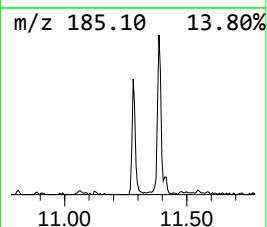
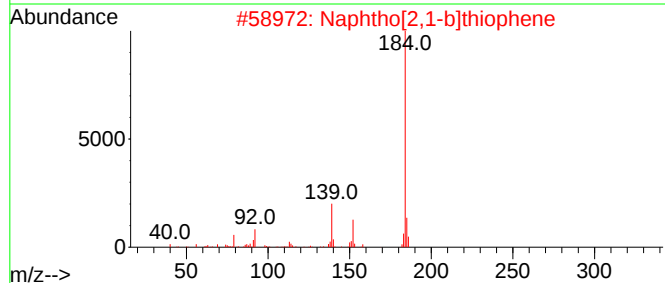
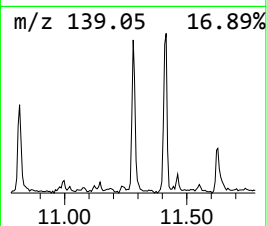
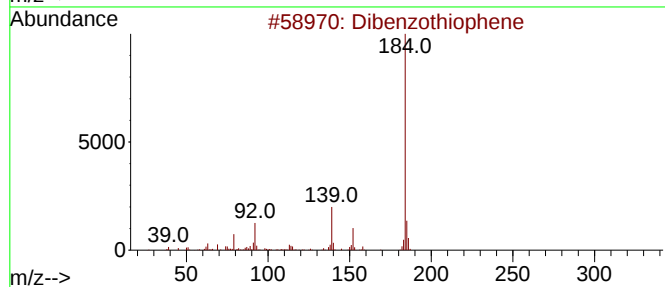
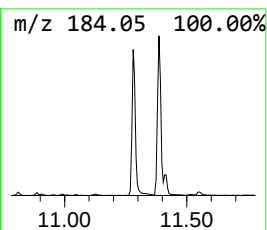
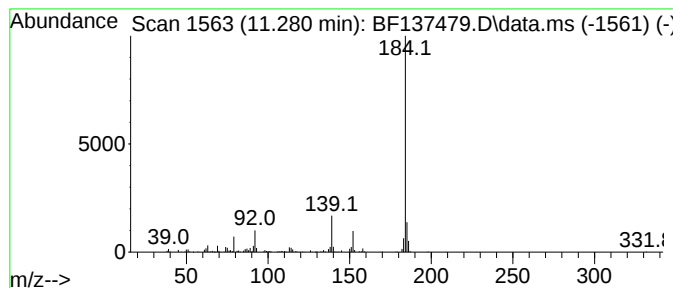
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Dibenzothiophene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.280	2.74 ng	295021	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95
3			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	94
4			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	93
5			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91



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Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
Operator : CG\JU
Sample : P1662-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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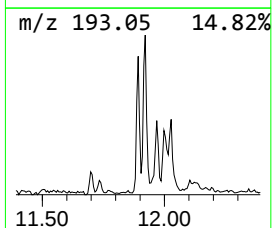
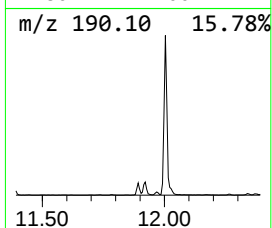
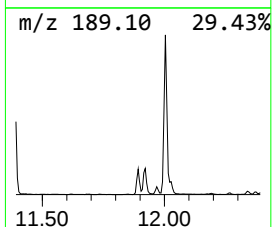
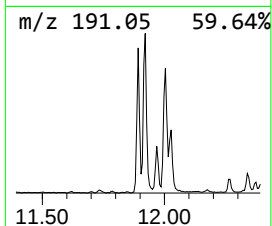
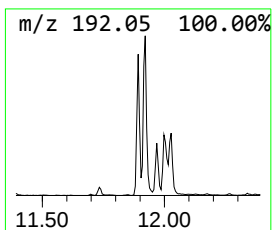
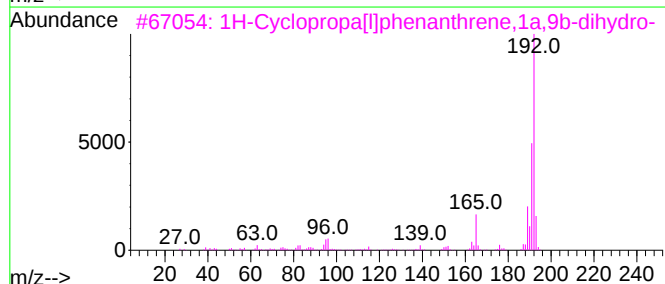
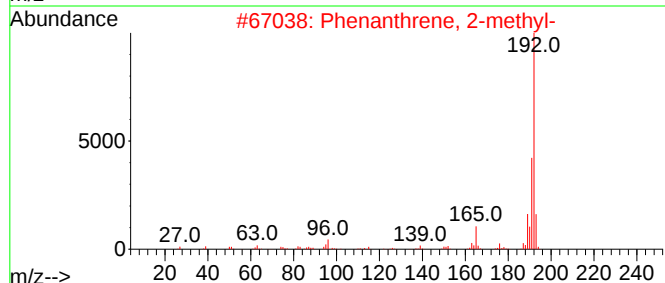
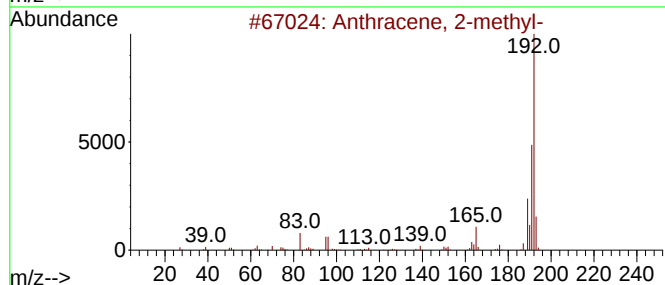
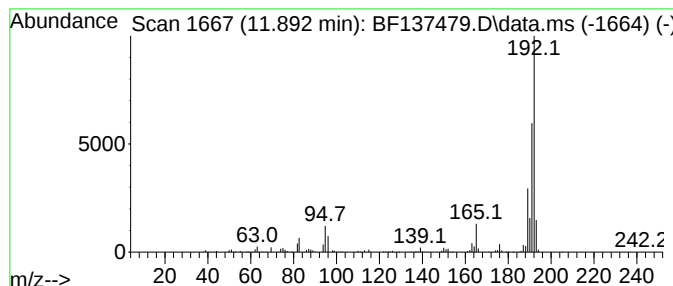
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.892	2.83 ng	303900	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
Operator : CG\JU
Sample : P1662-01
Misc :
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Instrument :
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ClientSampleId :
ETGI-352

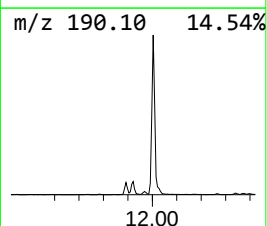
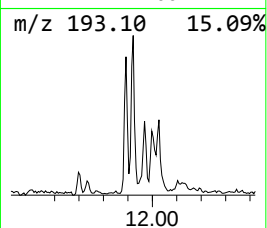
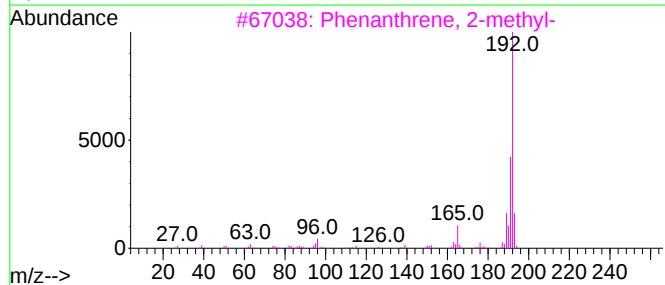
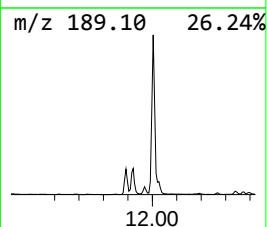
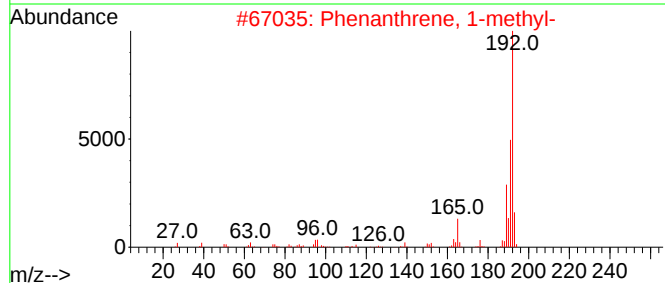
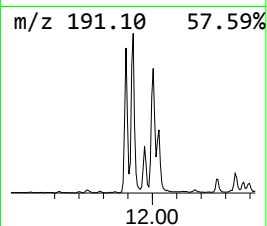
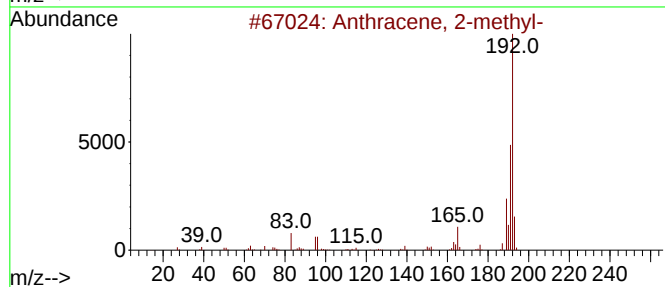
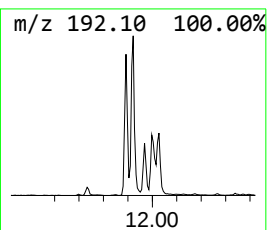
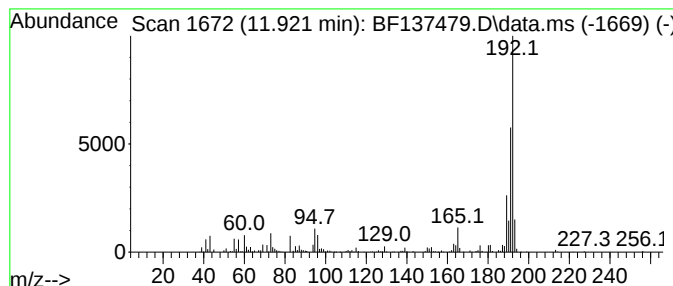
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.921	4.91 ng	528046	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
Operator : CG\JU
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ClientSampleId :
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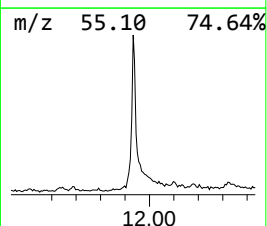
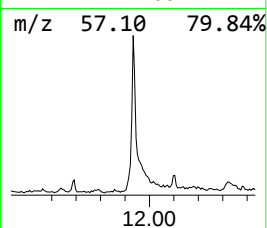
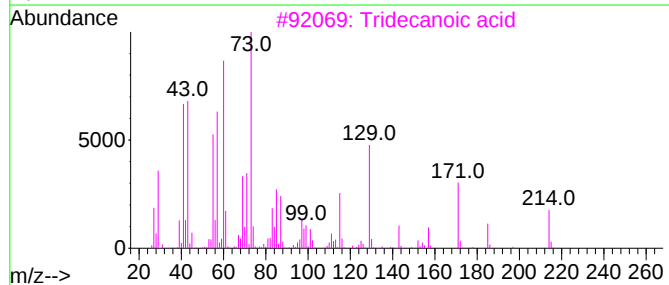
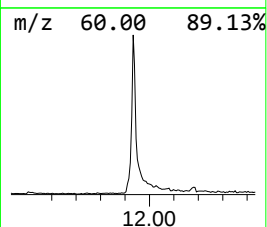
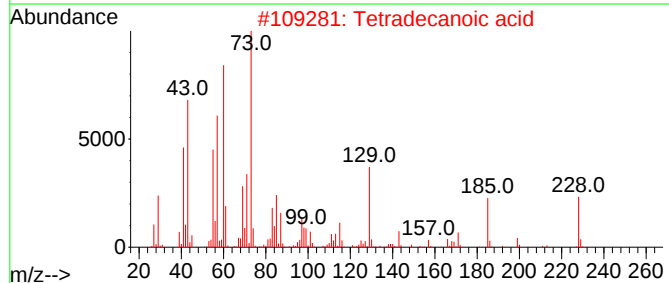
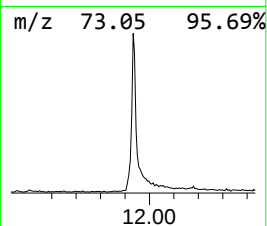
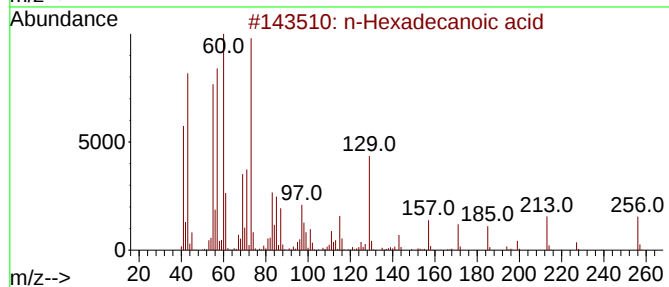
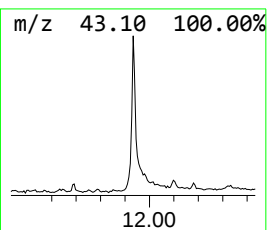
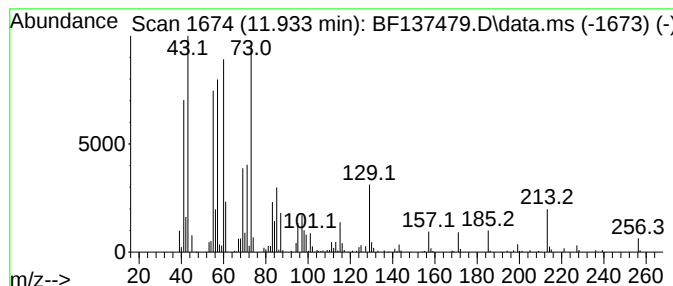
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.933	5.52 ng	594019	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3			Tridecanoic acid	214	C13H26O2	000638-53-9	74
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	72
5			Dodecanoic acid	200	C12H24O2	000143-07-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
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Sample : P1662-01
Misc :
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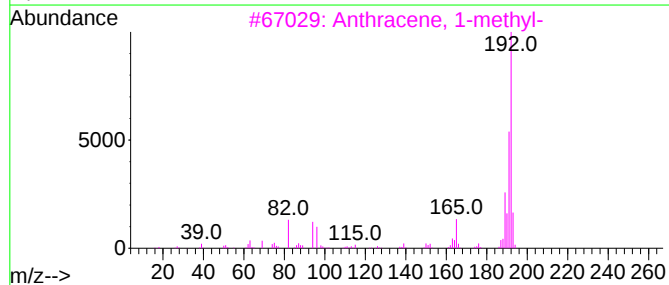
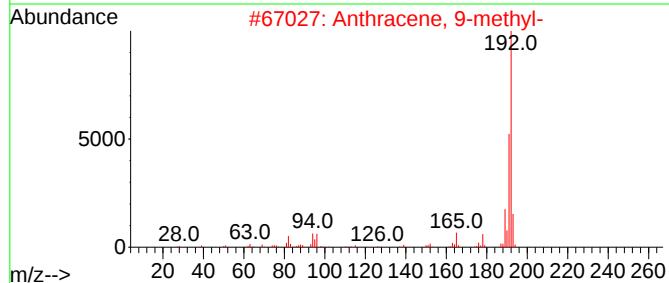
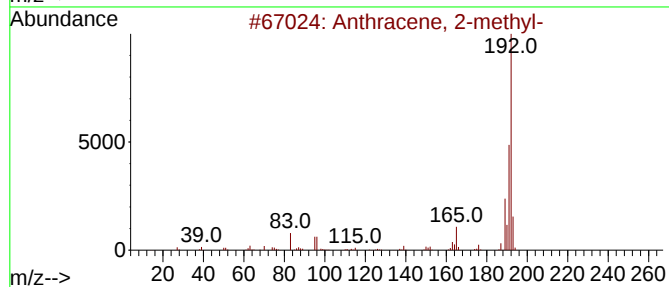
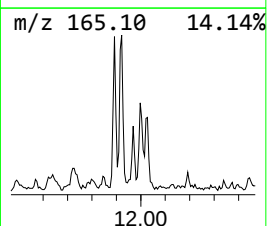
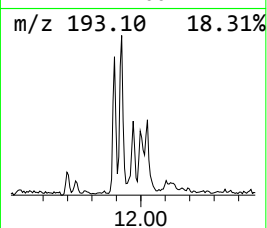
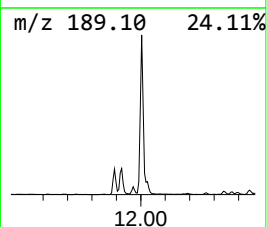
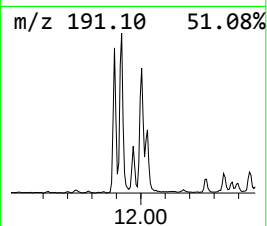
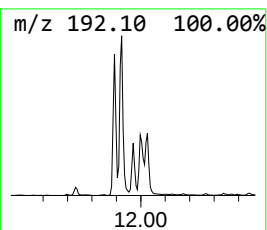
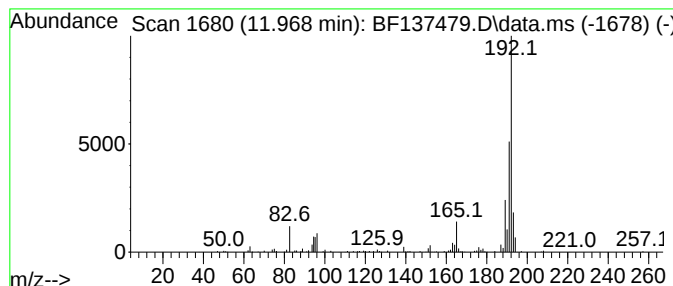
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Anthracene, 9-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.969	2.14 ng	229757	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
5			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
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Sample : P1662-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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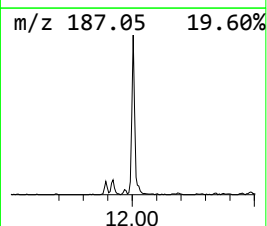
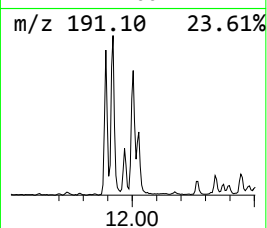
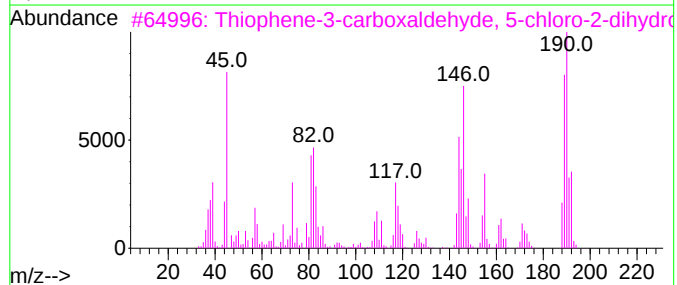
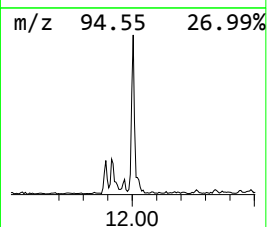
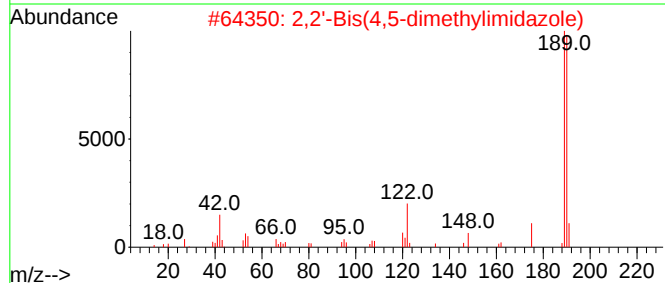
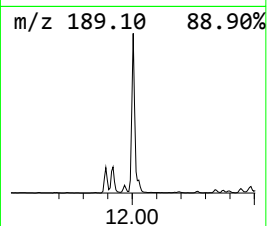
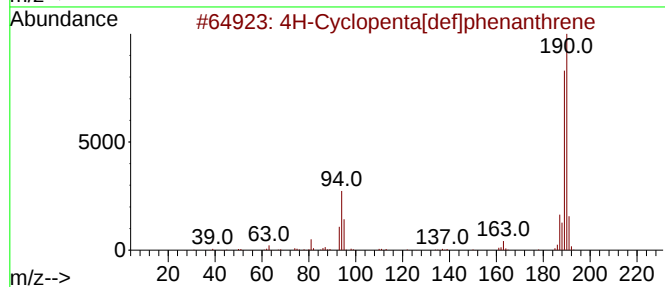
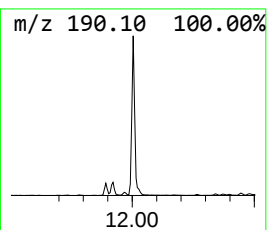
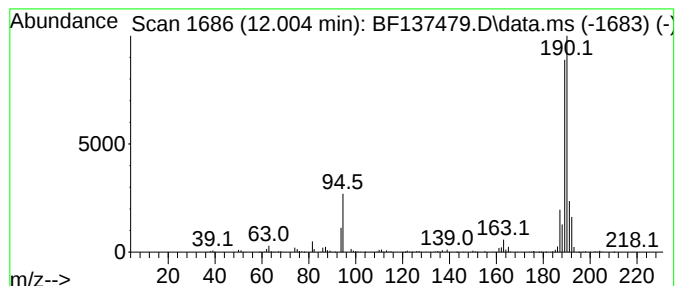
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.004	8.13 ng	875033	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
3			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	38
4			6,7-Methylenedioxy-4(3H)-quinazo...	190	C9H6N2O3	028310-11-4	33
5			1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
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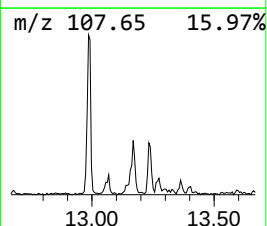
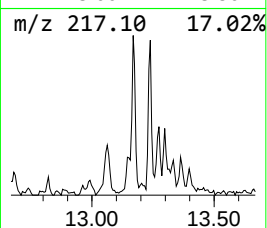
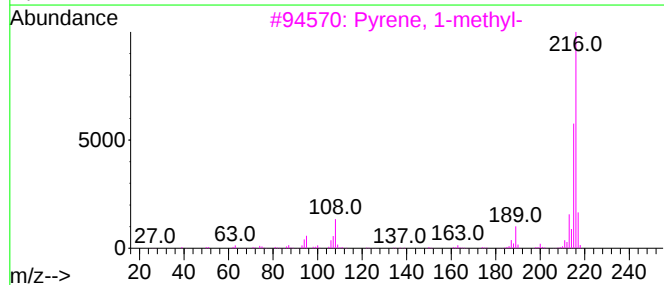
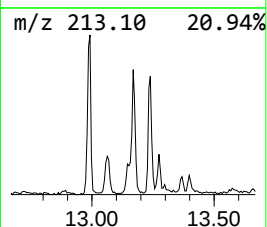
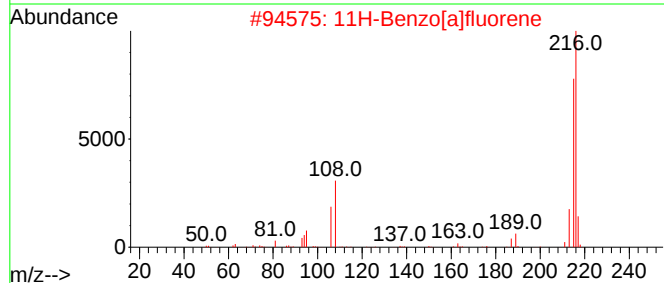
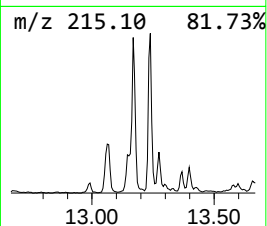
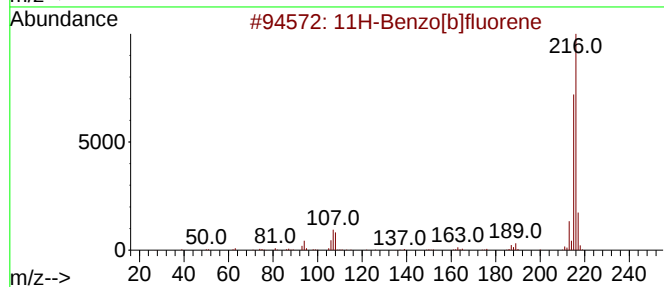
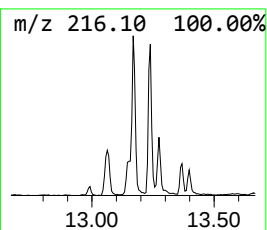
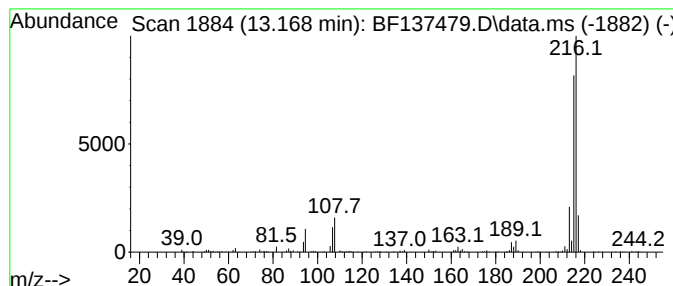
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.168	2.39 ng	247053	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	80
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	70



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Data File : BF137479.D
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Operator : CG\JU
Sample : P1662-01
Misc :
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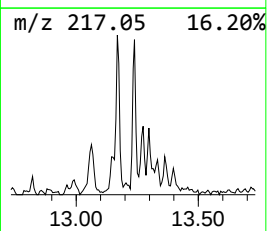
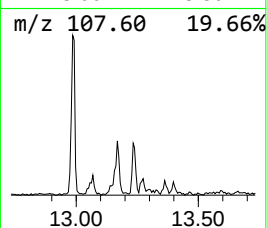
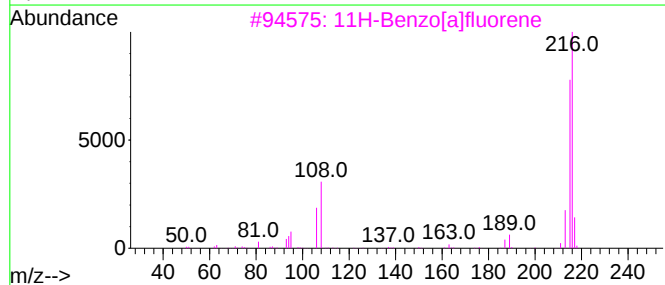
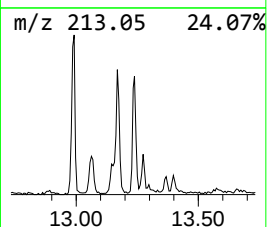
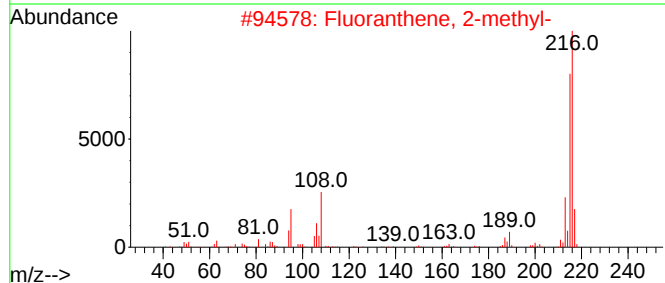
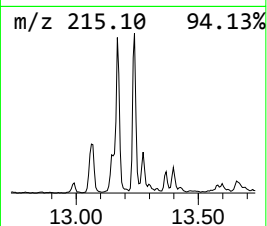
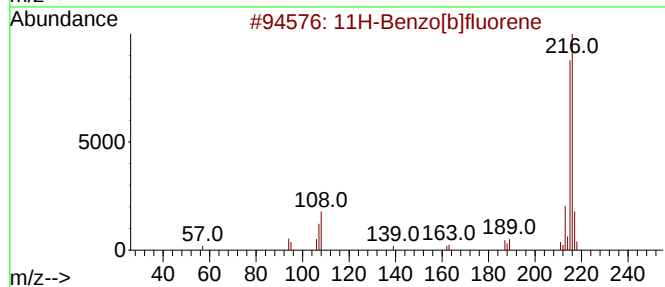
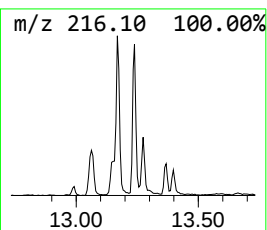
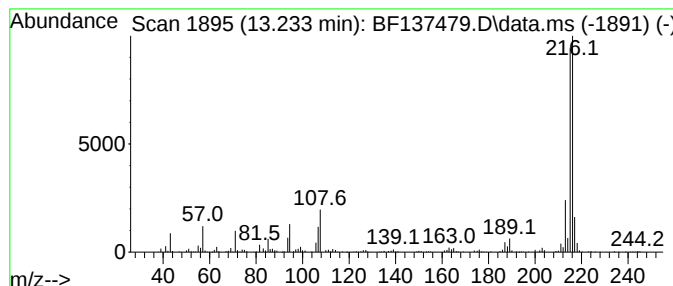
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Fluoranthene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.233	2.98 ng	307214	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



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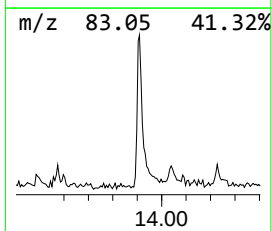
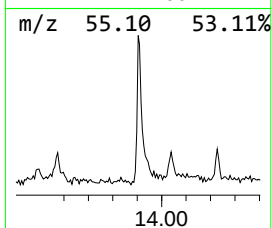
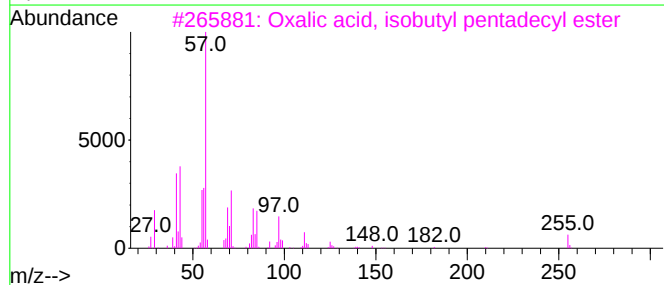
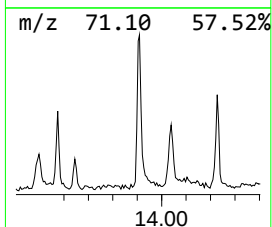
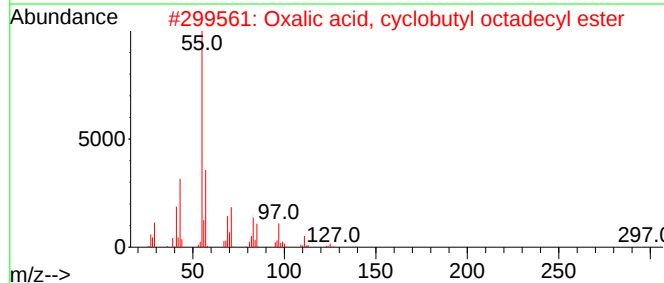
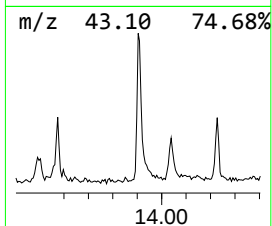
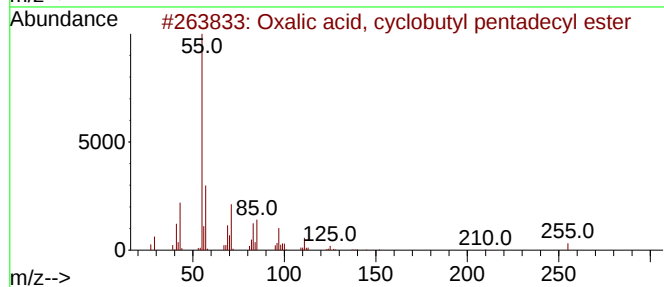
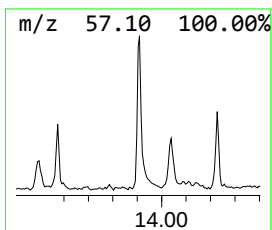
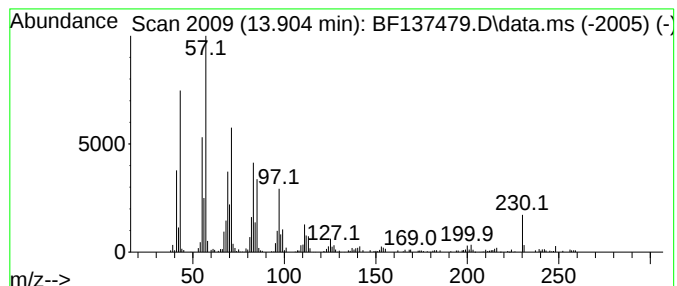
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Oxalic acid, cyclobutyl pen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	2.81 ng	289572	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclobutyl pentadec...	354	C21H38O4	1010309-70-5	91
2			Oxalic acid, cyclobutyl octadecy...	396	C24H44O4	1000309-70-8	91
3			Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	91
4			Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	90
5			Pentadecafluorooctanoic acid, pe...	624	C23H31F15O2	1000406-04-5	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
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Sample : P1662-01
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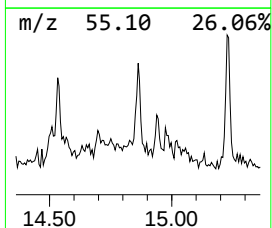
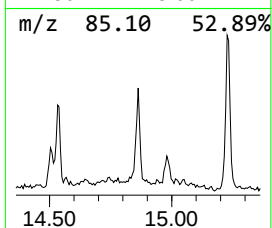
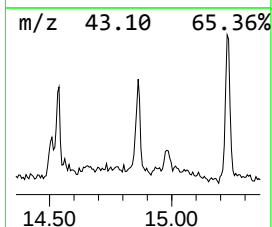
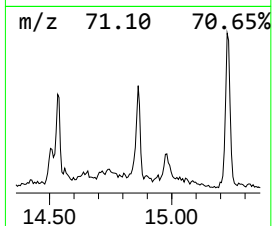
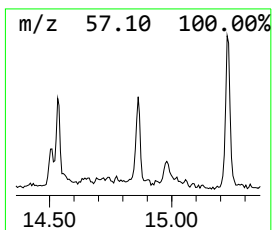
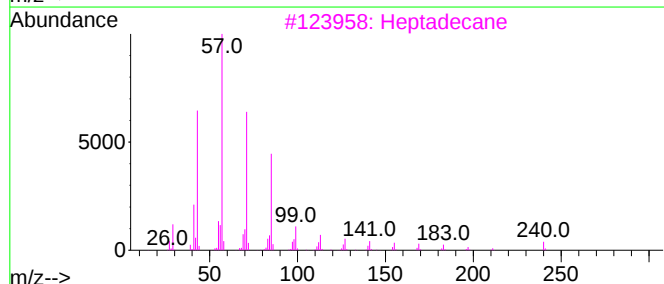
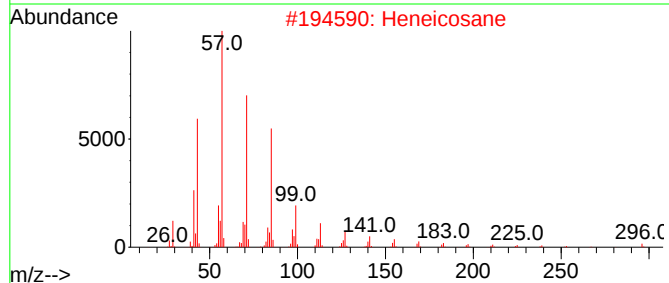
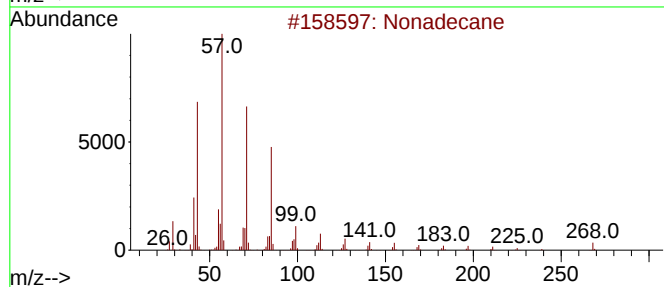
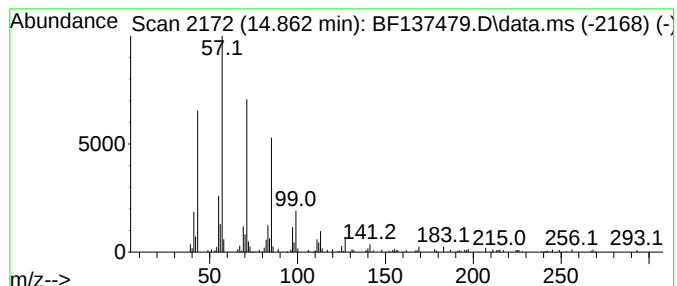
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.862	2.09 ng	130357	Perylene-d12	15.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	94
2			Heneicosane	296	C21H44	000629-94-7	91
3			Heptadecane	240	C17H36	000629-78-7	87
4			Octadecane	254	C18H38	000593-45-3	86
5			Hexadecane	226	C16H34	000544-76-3	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
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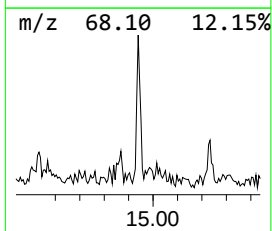
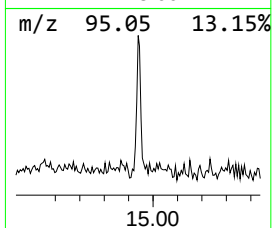
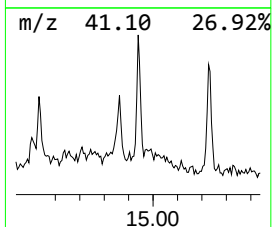
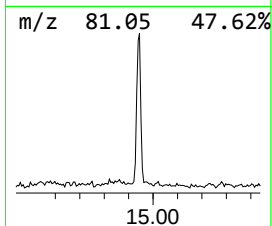
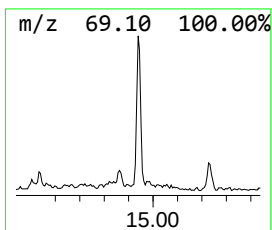
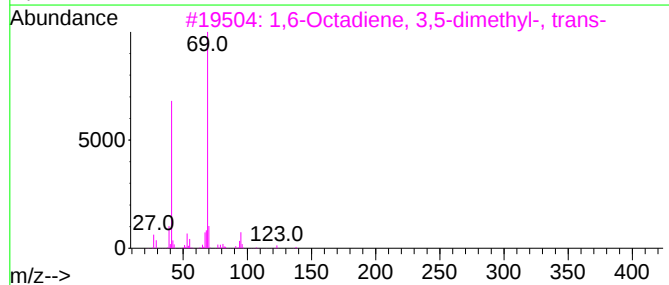
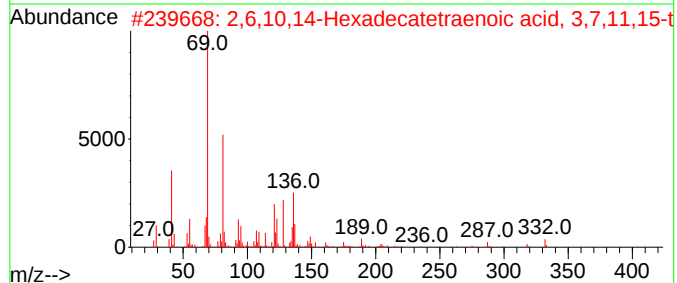
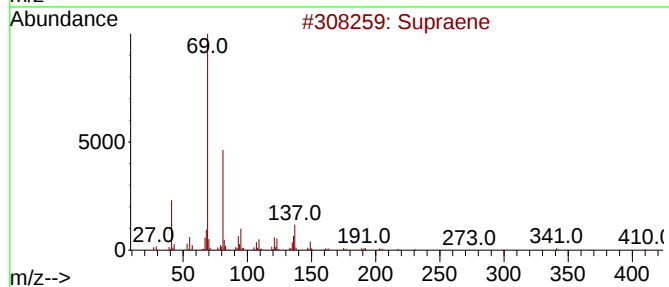
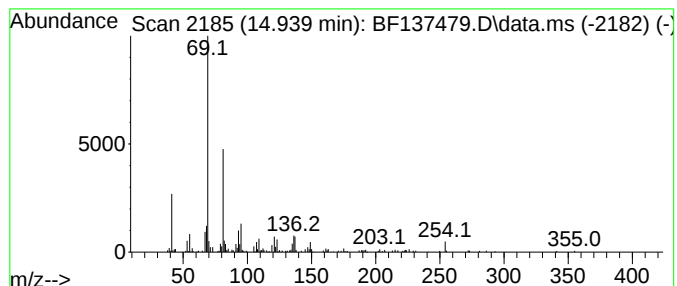
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Supraene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.939	2.10 ng	130789	Perylene-d12	15.574	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	87
2	2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	72
3	1,6-Octadiene, 3,5-dimethyl-, tr...	138	C10H18	074630-87-8	53
4	1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	53
5	Trideca-1,7,11-triene-1,1-dicarb...	270	C18H26N2	1000161-46-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
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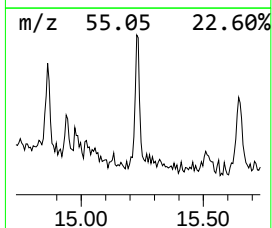
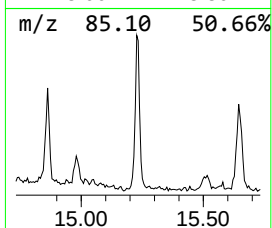
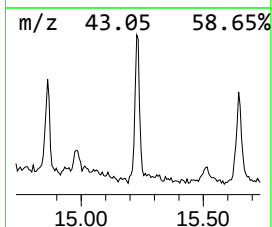
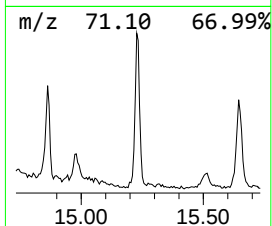
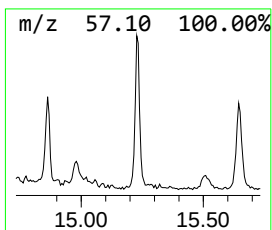
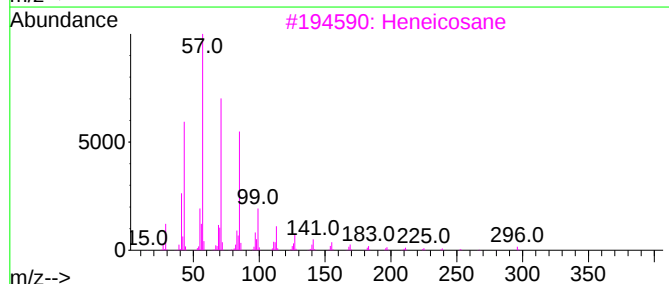
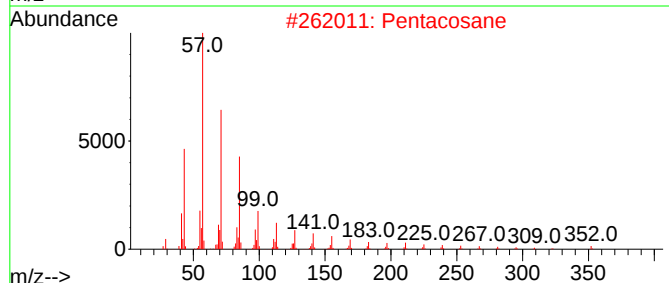
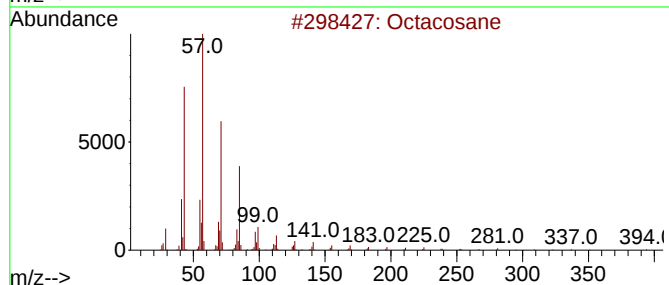
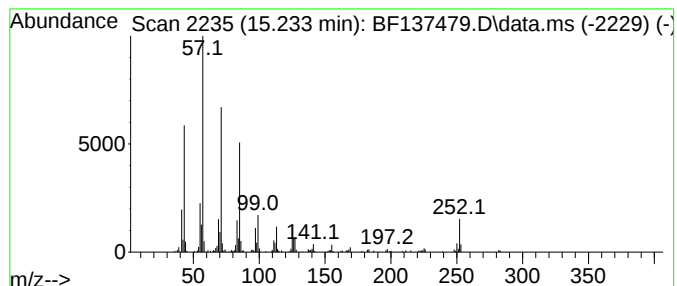
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Octacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.233	3.71 ng	231158	Perylene-d12	15.574

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane	394	C28H58	000630-02-4	87
2		Pentacosane	352	C25H52	000629-99-2	80
3		Heneicosane	296	C21H44	000629-94-7	76
4		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	74
5		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
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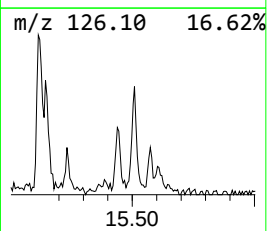
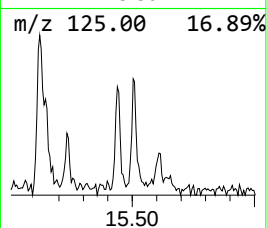
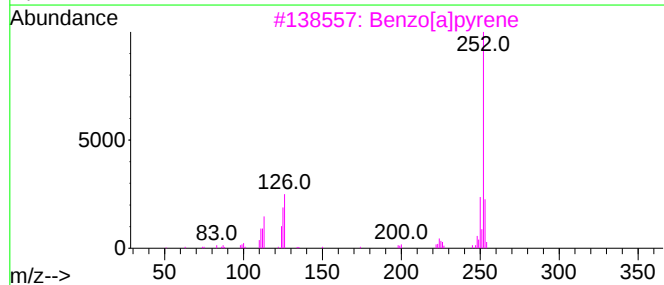
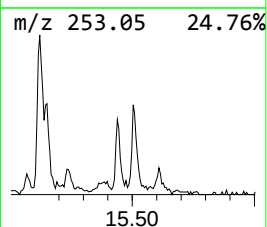
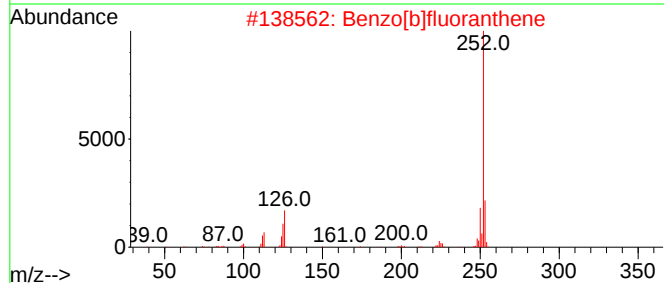
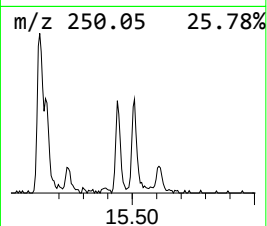
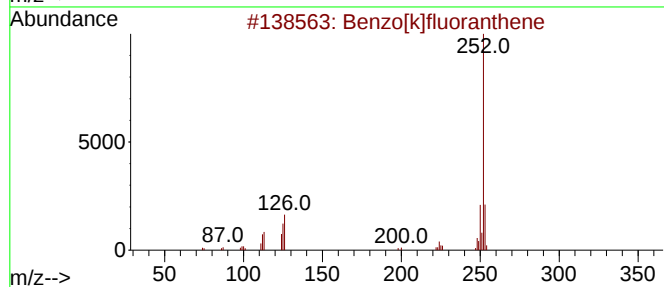
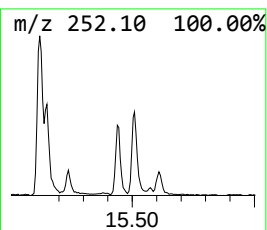
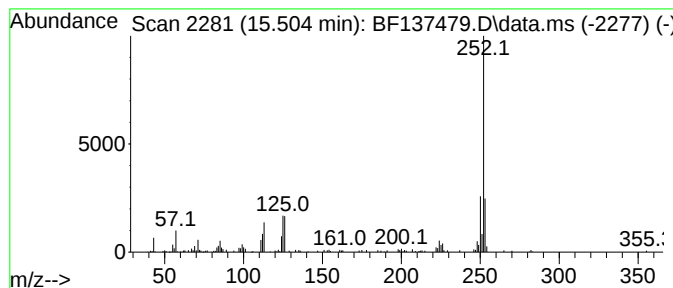
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Benzo[j]fluoranthene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.504	2.31 ng	144098	Perylene-d12	15.574

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
2		Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
3		Benzo[a]pyrene	252	C20H12	000050-32-8	95
4		Benzo[j]fluoranthene	252	C20H12	000205-82-3	93
5		Perylene	252	C20H12	000198-55-0	93



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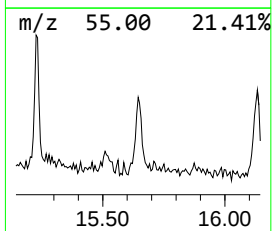
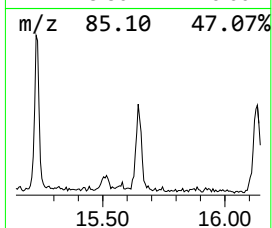
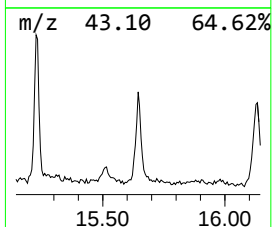
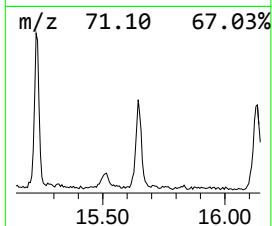
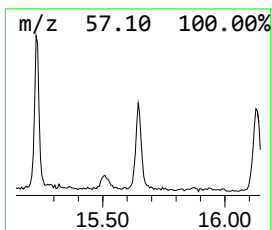
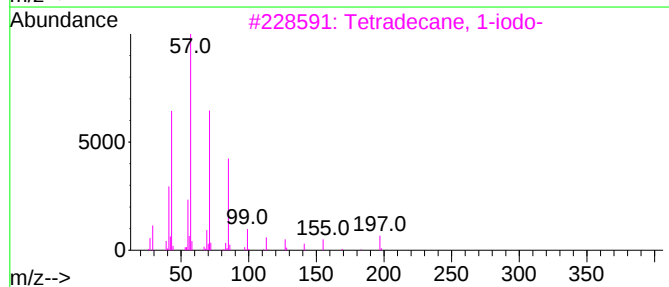
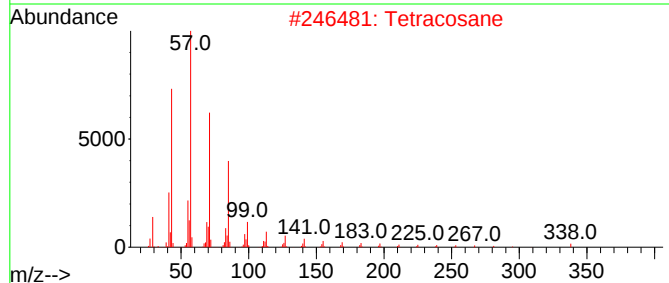
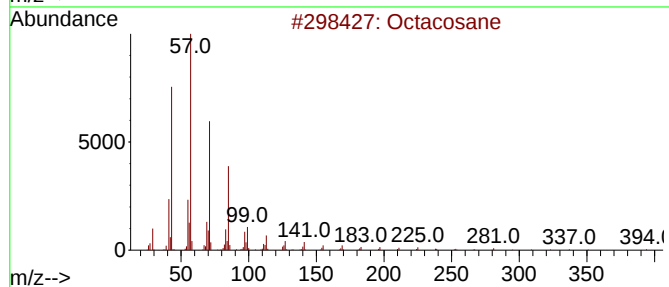
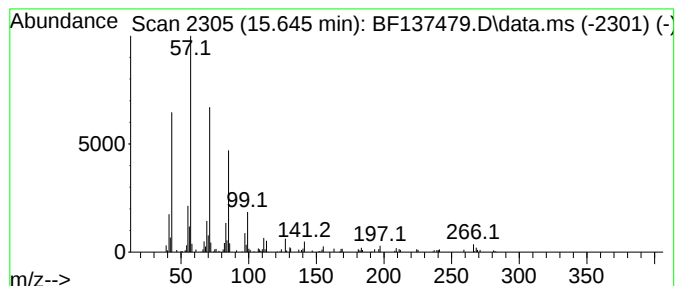
Instrument :
BNA_F
ClientSampleId :
ETGI-352

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Tetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.645	2.74 ng	170710	Perylene-d12	15.574	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	83
2	Tetracosane	338	C24H50	000646-31-1	83
3	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	80
4	Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	80
5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
Operator : CG\JU
Sample : P1662-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-352

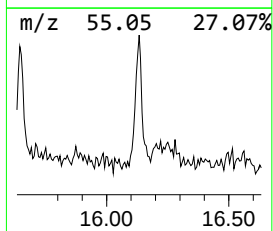
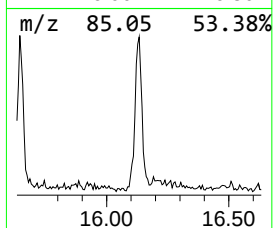
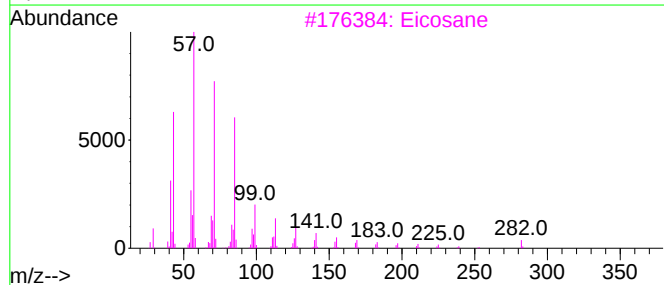
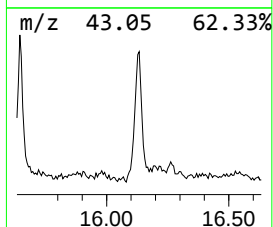
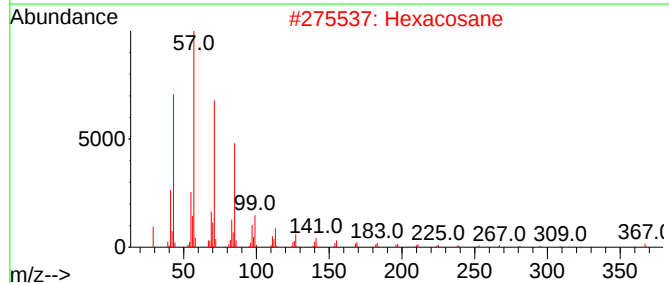
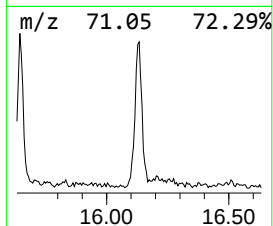
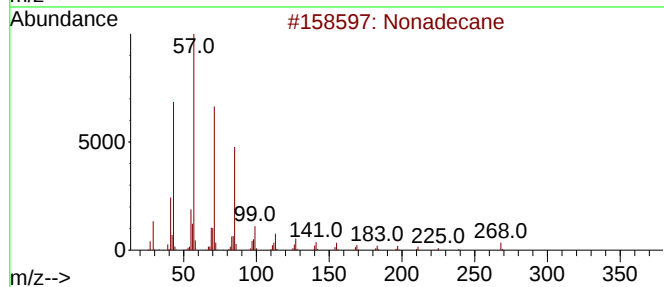
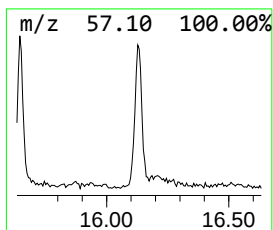
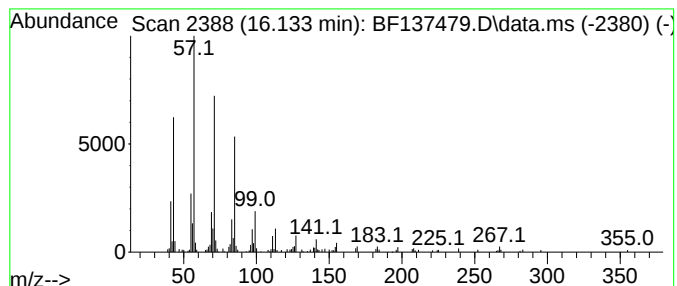
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.133	3.24 ng	202221	Perylene-d12	15.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			Hexacosane	366	C26H54	000630-01-3	91
3			Eicosane	282	C20H42	000112-95-8	91
4			Tetracosane	338	C24H50	000646-31-1	91
5			Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030424\
Data File : BF137479.D
Acq On : 04 Mar 2024 14:20
Operator : CG\JU
Sample : P1662-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-352

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.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...	2.187	29.3	ng	1918680	1	6.857	1310500	20.0
Dibenzofuran, 4...	10.598	2.3	ng	244862	3	9.892	2136570	20.0
9H-Fluorene, 1-...	10.992	2.2	ng	239576	4	11.386	2151310	20.0
Dibenzothiophene	11.280	2.7	ng	295021	4	11.386	2151310	20.0
Anthracene, 2-m...	11.892	2.8	ng	303900	4	11.386	2151310	20.0
Phenanthrene, 1...	11.921	4.9	ng	528046	4	11.386	2151310	20.0
n-Hexadecanoic ...	11.933	5.5	ng	594019	4	11.386	2151310	20.0
Anthracene, 9-m...	11.969	2.1	ng	229757	4	11.386	2151310	20.0
4H-Cyclopenta[d...	12.004	8.1	ng	875033	4	11.386	2151310	20.0
11H-Benzo[b]flu...	13.168	2.4	ng	247053	5	14.045	2063140	20.0
Fluoranthene, 2...	13.233	3.0	ng	307214	5	14.045	2063140	20.0
Oxalic acid, cy...	13.904	2.8	ng	289572	5	14.045	2063140	20.0
Nonadecane	14.862	2.1	ng	130357	6	15.574	1247050	20.0
Supraene	14.939	2.1	ng	130789	6	15.574	1247050	20.0
Octacosane	15.233	3.7	ng	231158	6	15.574	1247050	20.0
Benzo[j]fluoran...	15.504	2.3	ng	144098	6	15.574	1247050	20.0
Tetracosane	15.645	2.7	ng	170710	6	15.574	1247050	20.0
Hexacosane	16.133	3.2	ng	202221	6	15.574	1247050	20.0