

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081723\
 Data File : BF134828.D
 Acq On : 17 Aug 2023 21:05
 Operator : CG\JU
 Sample : 04042-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-376

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-BF080123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF134828.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.422	563	567	571	rBV	2507778	2483148	94.11%	8.792%
2	6.428	733	738	741	rBV	2647439	2471380	93.67%	8.750%
3	6.792	796	800	803	rBV	1127964	852410	32.31%	3.018%
4	7.351	890	895	898	rBV	1809923	1562780	59.23%	5.533%
5	8.069	1013	1017	1020	rBV	1330068	1089921	41.31%	3.859%
6	9.145	1196	1200	1204	rBV	2946647	2638519	100.00%	9.342%
7	9.263	1218	1220	1224	rVB	32719	29225	1.11%	0.103%
8	9.686	1288	1292	1297	rBV	169641	136711	5.18%	0.484%
9	9.827	1312	1316	1319	rBV	1249829	1090296	41.32%	3.860%
10	10.027	1346	1350	1354	rBV	37181	37059	1.40%	0.131%
11	10.374	1406	1409	1415	rVB2	38696	50756	1.92%	0.180%
12	10.616	1446	1450	1453	rBV	1632140	1417435	53.72%	5.019%
13	10.739	1468	1471	1479	rBV4	29813	31691	1.20%	0.112%
14	10.921	1497	1502	1505	rBV5	17237	28761	1.09%	0.102%
15	10.980	1509	1512	1515	rBV3	30930	29257	1.11%	0.104%
16	11.210	1549	1551	1556	rVB	30015	30102	1.14%	0.107%
17	11.316	1565	1569	1571	rBV	1091608	931974	35.32%	3.300%
18	11.339	1571	1573	1576	rVV	391379	309613	11.73%	1.096%
19	11.386	1578	1581	1584	rVB	102370	94482	3.58%	0.335%
20	11.421	1584	1587	1591	rBV2	37917	41345	1.57%	0.146%
21	11.774	1643	1647	1650	rBV4	33094	39517	1.50%	0.140%
22	11.816	1651	1654	1657	rVV	76502	87157	3.30%	0.309%
23	11.851	1657	1660	1665	rVV2	464357	452565	17.15%	1.602%
24	11.892	1665	1667	1670	rVV	53046	58518	2.22%	0.207%
25	11.927	1670	1673	1676	rVV	137807	155079	5.88%	0.549%
26	11.951	1676	1677	1682	rVB	69152	48769	1.85%	0.173%
27	12.116	1702	1705	1707	rBV2	45028	34204	1.30%	0.121%
28	12.139	1707	1709	1716	rVB2	41052	40141	1.52%	0.142%
29	12.368	1742	1748	1751	rBV	86245	98818	3.75%	0.350%
30	12.404	1751	1754	1756	rVV2	45379	56884	2.16%	0.201%
31	12.457	1760	1763	1767	rVV3	45289	56859	2.15%	0.201%
32	12.533	1772	1776	1779	rBV	600122	561222	21.27%	1.987%
33	12.639	1789	1794	1797	rBV2	79007	132008	5.00%	0.467%
34	12.757	1808	1814	1818	rBV	572475	628420	23.82%	2.225%
35	12.821	1821	1825	1828	rVB2	31373	47895	1.82%	0.170%
36	12.910	1836	1840	1843	rBV	2203087	1873565	71.01%	6.634%

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37	12.980	1847	1852	1855	rBV3	66109	98986	3.75%	0.350%
38	13.068	1863	1867	1868	rBV4	56969	65357	2.48%	0.231%
39	13.092	1868	1871	1874	rVB2	116695	114207	4.33%	0.404%
40	13.157	1879	1882	1885	rVB2	79011	68163	2.58%	0.241%
41	13.192	1885	1888	1892	rBV	86560	85787	3.25%	0.304%
42	13.263	1896	1900	1902	rBV2	31533	42453	1.61%	0.150%
43	13.286	1902	1904	1907	rVB	68616	51042	1.93%	0.181%
44	13.315	1907	1909	1913	rBV	59724	53847	2.04%	0.191%
45	13.398	1917	1923	1932	rVV	771642	732284	27.75%	2.593%
46	13.598	1954	1957	1963	rVB	62532	87991	3.33%	0.312%
47	13.710	1972	1976	1979	rBV3	76667	110817	4.20%	0.392%
48	13.739	1979	1981	1983	rVV	67632	64620	2.45%	0.229%
49	13.762	1983	1985	1989	rVV2	69284	94002	3.56%	0.333%
50	13.821	1990	1995	2003	rVB3	131078	229772	8.71%	0.814%
51	13.962	2014	2019	2022	rBV2	1153830	1379387	52.28%	4.884%
52	13.992	2022	2024	2030	rVB	386492	337472	12.79%	1.195%
53	14.068	2035	2037	2039	rVB	45886	30814	1.17%	0.109%
54	14.315	2077	2079	2081	rBV2	36516	34575	1.31%	0.122%
55	14.351	2081	2085	2088	rVB	110986	129324	4.90%	0.458%
56	14.386	2088	2091	2095	rVB	74476	65975	2.50%	0.234%
57	14.445	2099	2101	2104	rVB2	64343	58397	2.21%	0.207%
58	14.474	2104	2106	2108	rBV2	66188	57791	2.19%	0.205%
59	14.833	2163	2167	2173	rBV	628135	641528	24.31%	2.271%
60	14.915	2178	2181	2183	rBV3	41776	42057	1.59%	0.149%
61	15.004	2192	2196	2200	rBV	444429	730907	27.70%	2.588%
62	15.115	2211	2215	2219	rVB	134582	164308	6.23%	0.582%
63	15.309	2244	2248	2254	rVB	296411	379474	14.38%	1.344%
64	15.368	2254	2258	2265	rBV	415017	544176	20.62%	1.927%
65	15.439	2265	2270	2273	rVV	853911	1069012	40.52%	3.785%
66	15.468	2273	2275	2282	rVB2	151620	205328	7.78%	0.727%
67	16.915	2516	2521	2531	rVB2	217849	444104	16.83%	1.572%
68	17.368	2592	2598	2604	rBV	148079	300520	11.39%	1.064%

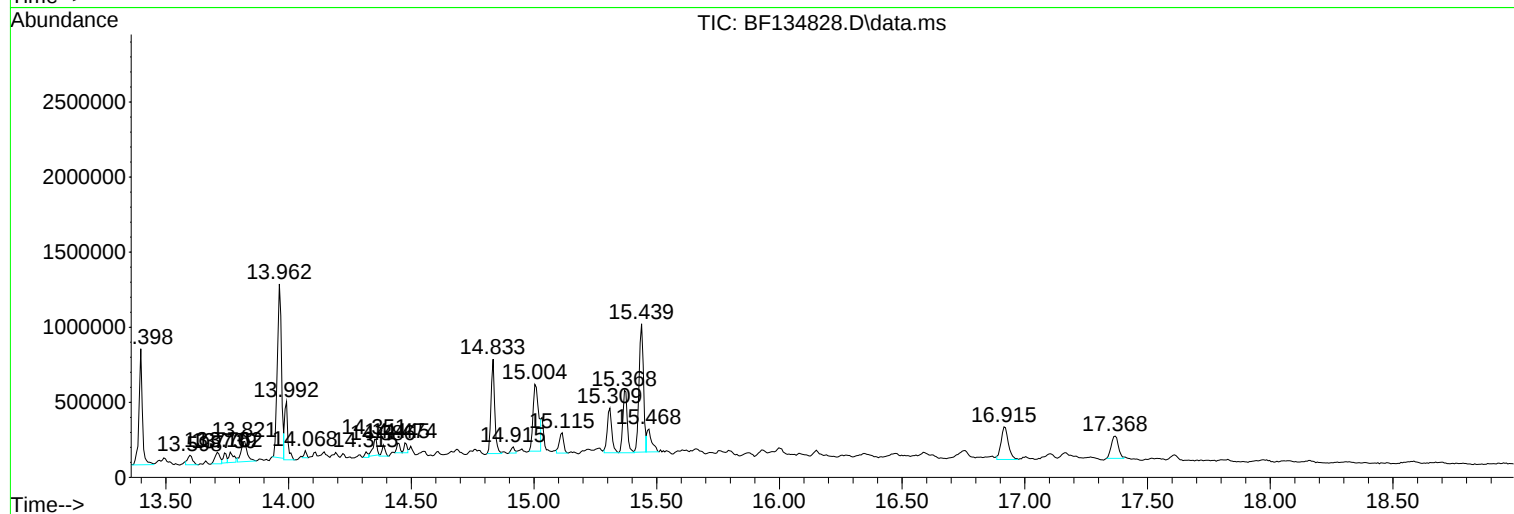
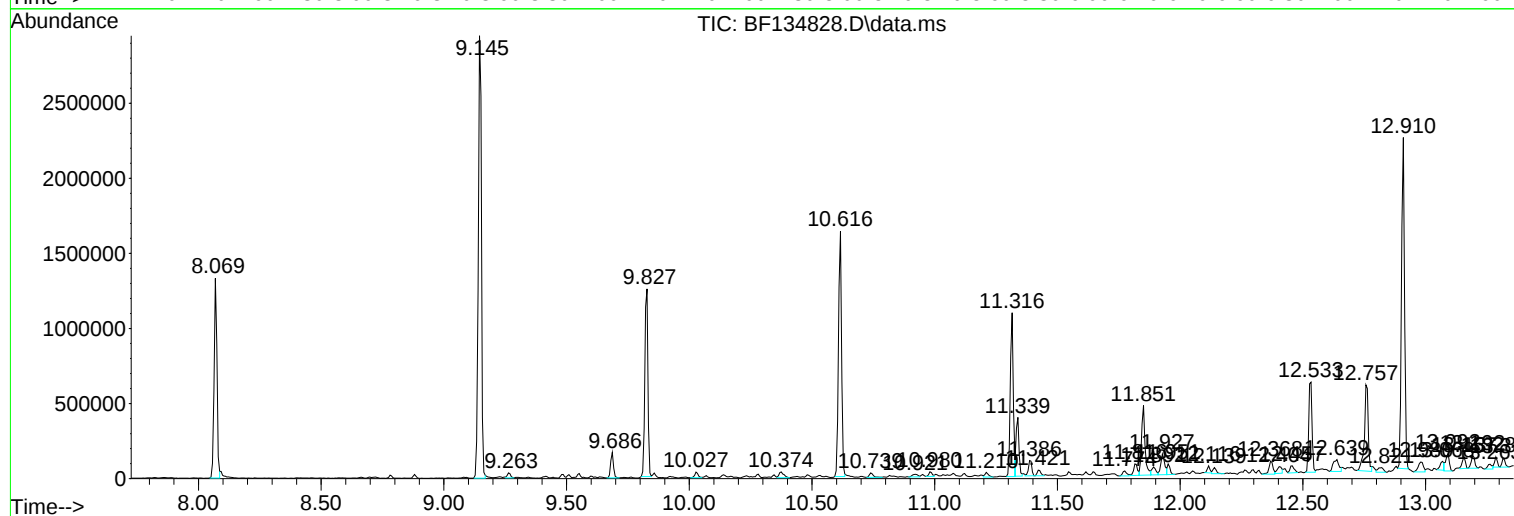
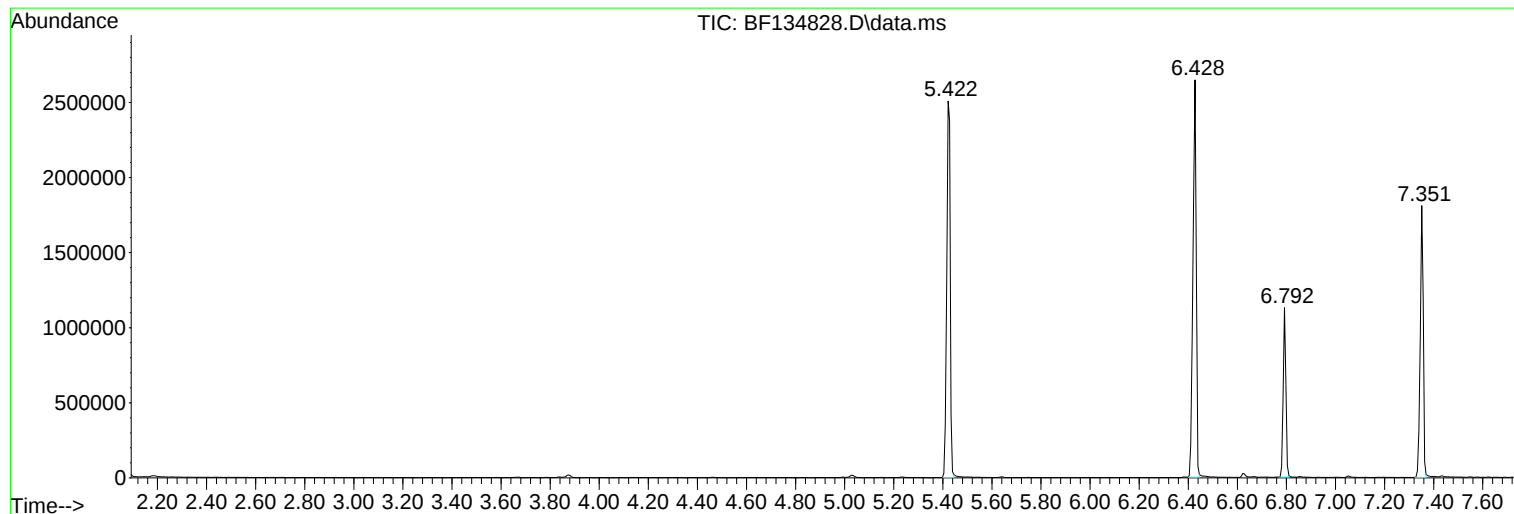
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BNA_F
ClientSampleId :
TP-376

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.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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Instrument :
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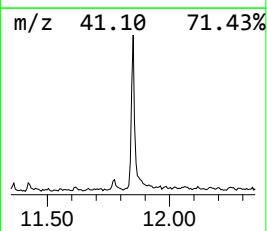
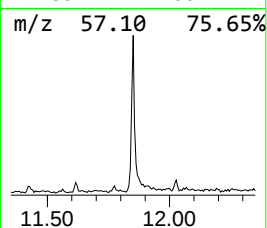
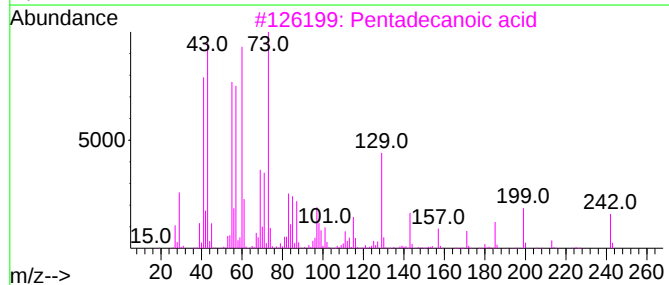
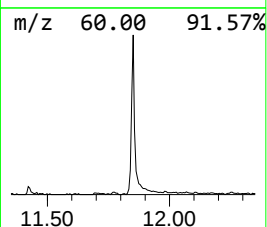
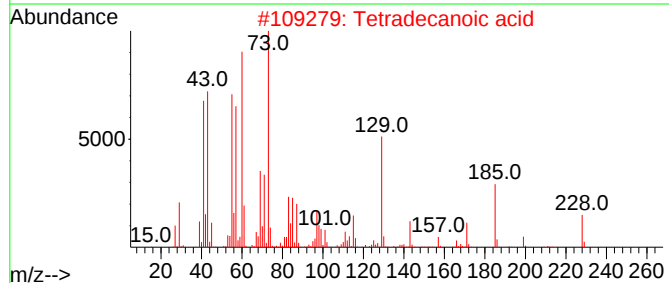
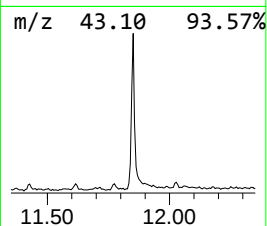
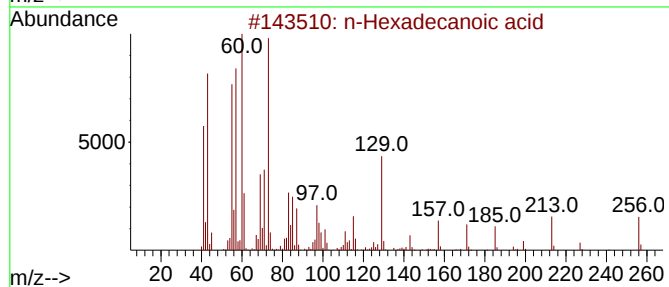
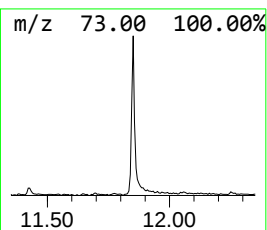
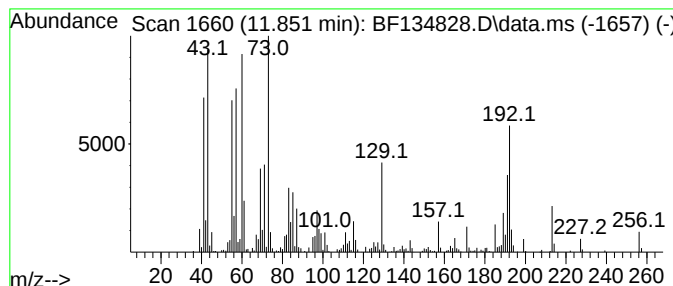
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.851	9.71 ng	452565	Phenanthrene-d10	11.316

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4		Tridecanoic acid	214	C13H26O2	000638-53-9	86
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



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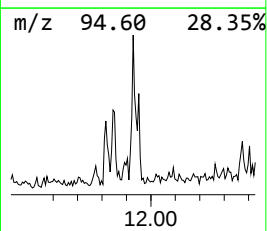
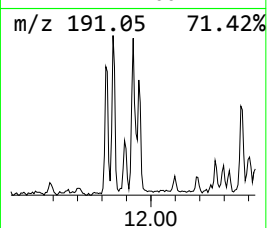
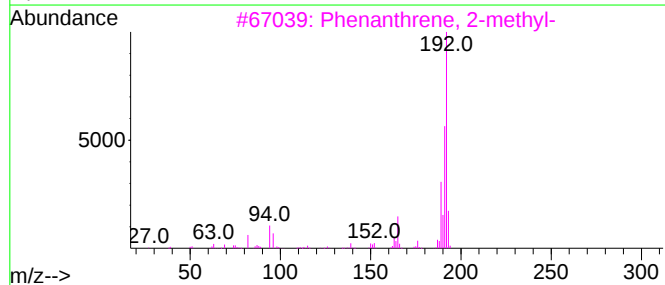
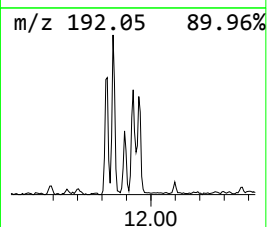
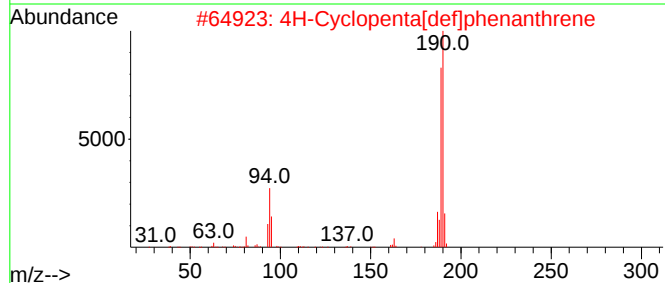
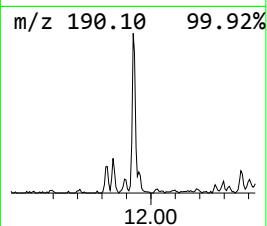
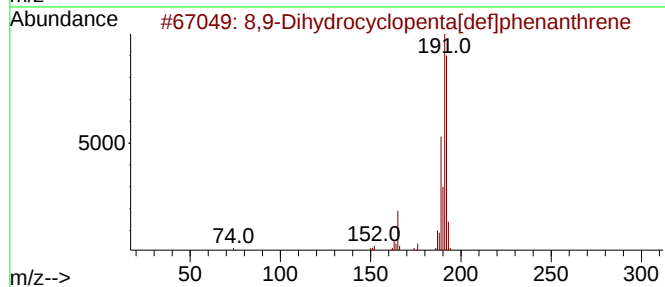
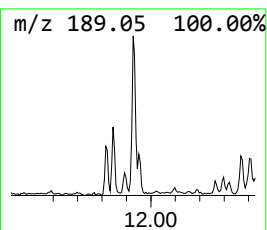
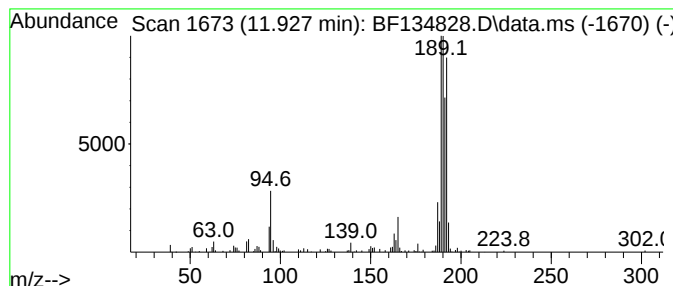
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 8,9-Dihydrocyclopenta[def]p... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
11.927	3.33 ng	155079	Phenanthrene-d10			11.316
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	8,9-Dihydrocyclopenta[def]phenan...		192	C15H12	027410-55-5	74
2	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	46
3	Phenanthrene, 2-methyl-		192	C15H12	002531-84-2	42
4	Phenanthrene, 4-methyl-		192	C15H12	000832-64-4	38
5	Anthracene, 1-methyl-		192	C15H12	000610-48-0	38



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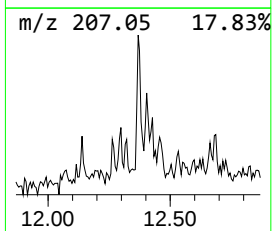
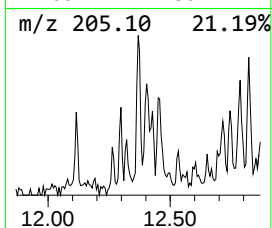
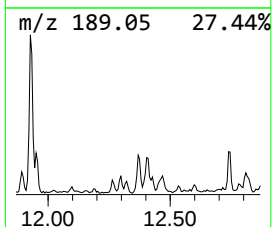
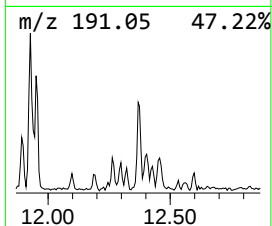
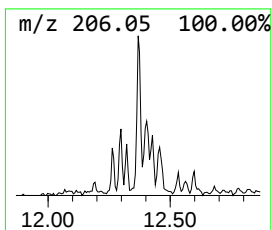
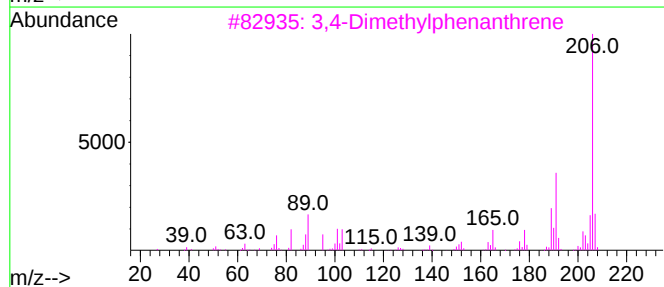
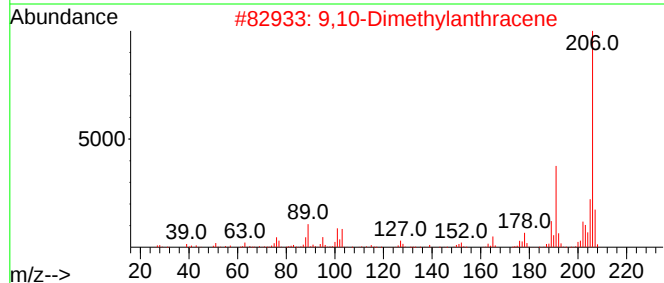
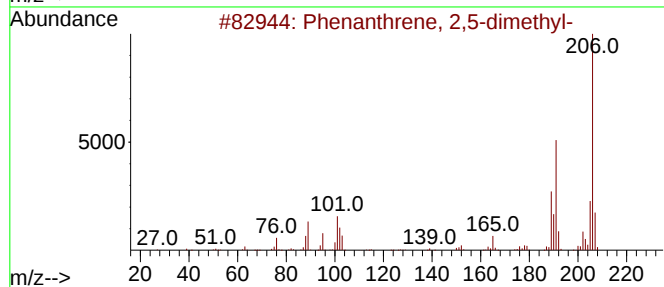
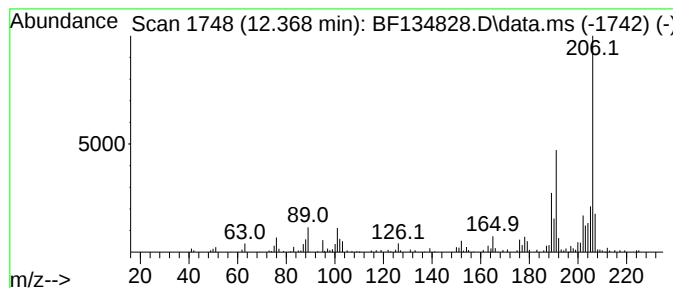
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Peak Number 3 Phenanthrene, 2,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.368	2.12 ng	98818	Phenanthrene-d10	11.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2			9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
3			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	93
4			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91
5			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90



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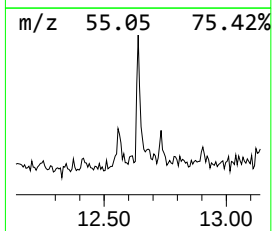
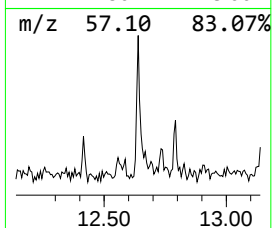
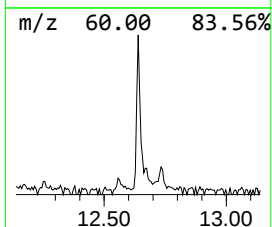
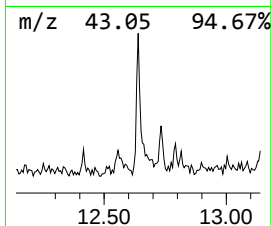
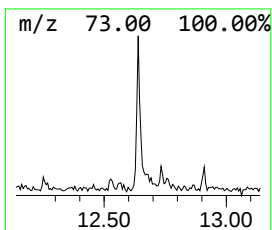
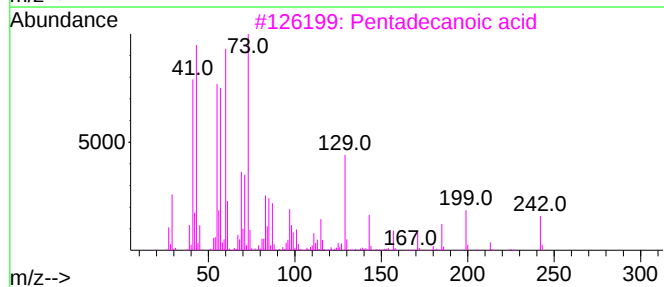
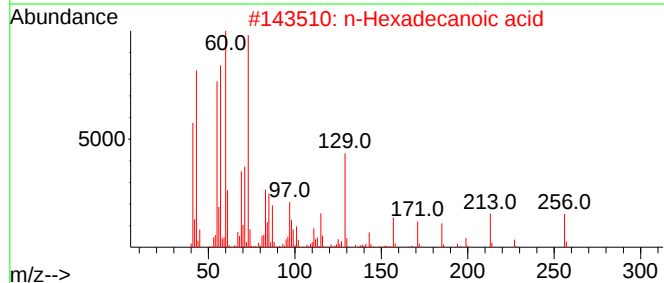
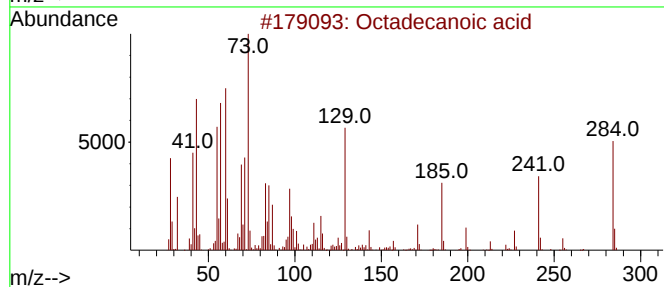
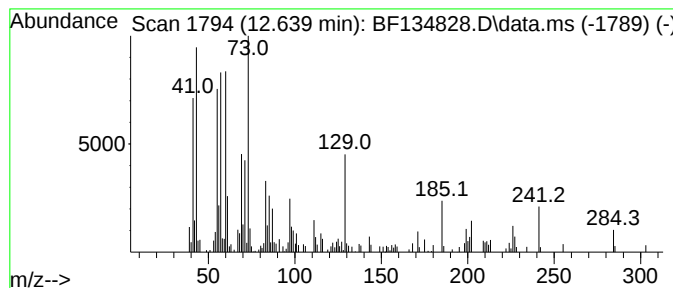
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.639	2.83 ng	132008	Phenanthrene-d10	11.316

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	87
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081723\
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Acq On : 17 Aug 2023 21:05
Operator : CG\JU
Sample : 04042-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-376

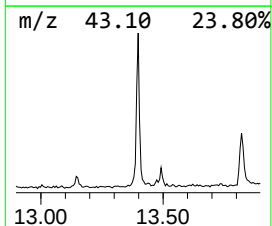
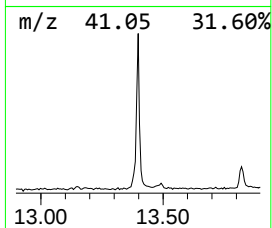
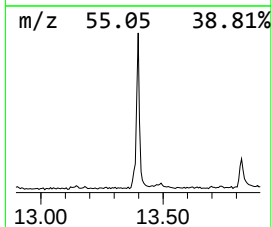
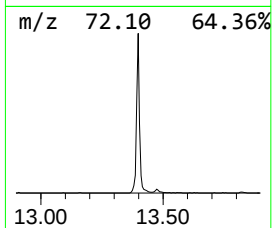
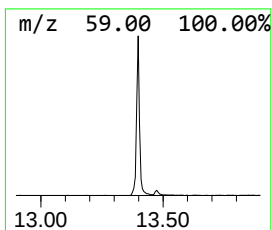
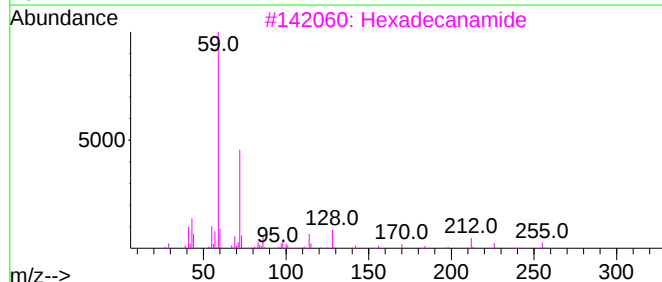
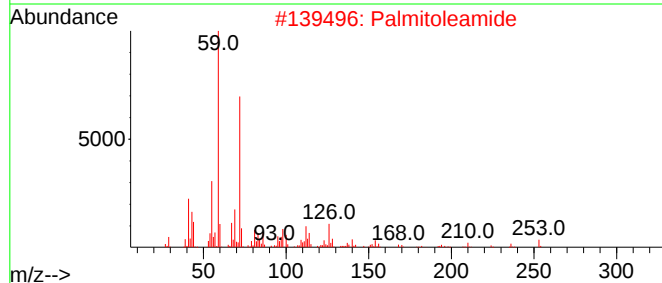
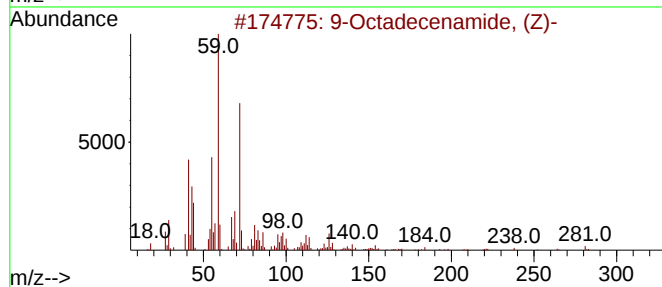
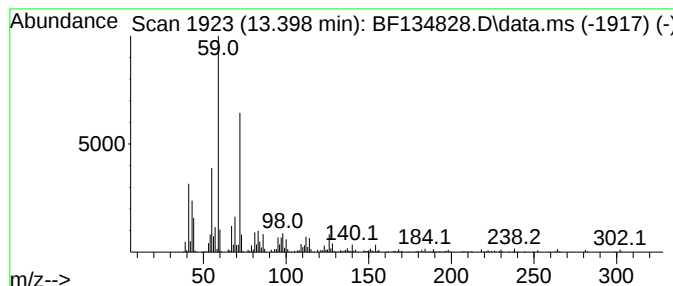
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 9-Octadecenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.398	10.62 ng	732284	Chrysene-d12	13.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	97
2			Palmitoleamide	253	C16H31NO	106010-22-4	86
3			Hexadecanamide	255	C16H33NO	000629-54-9	78
4			Nonanamide	157	C9H19NO	001120-07-6	59
5			Octadecanamide	283	C18H37NO	000124-26-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081723\
Data File : BF134828.D
Acq On : 17 Aug 2023 21:05
Operator : CG\JU
Sample : 04042-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-376

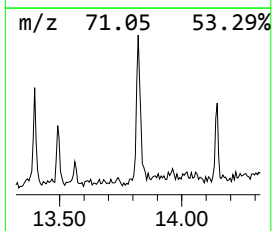
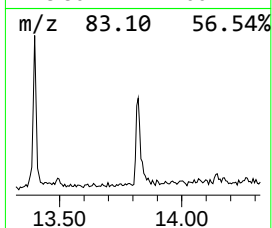
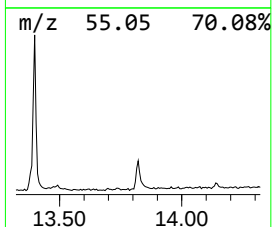
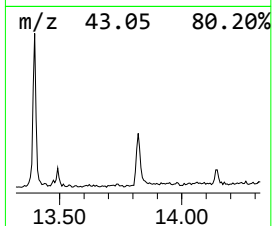
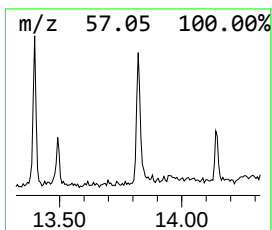
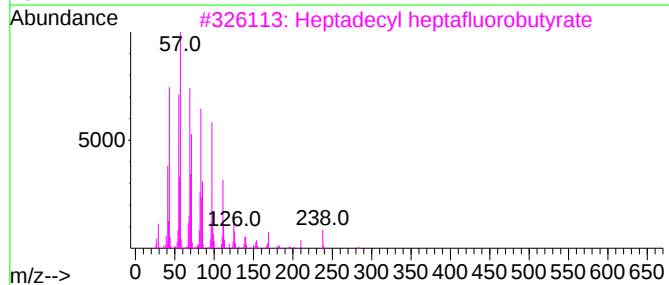
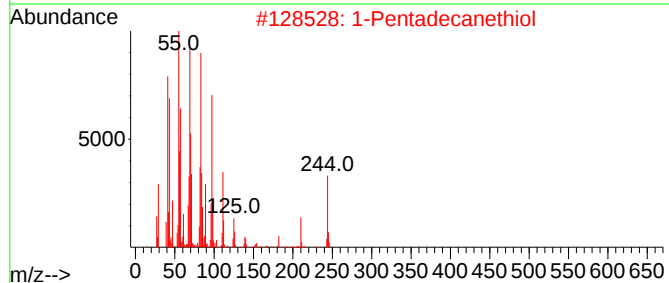
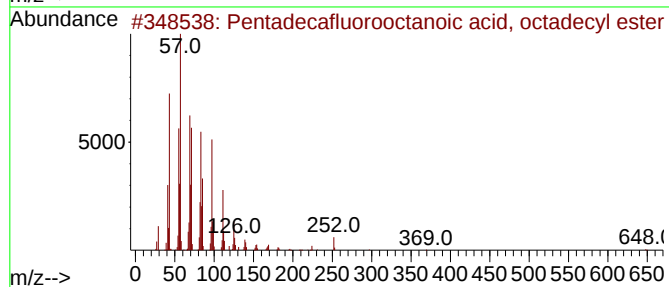
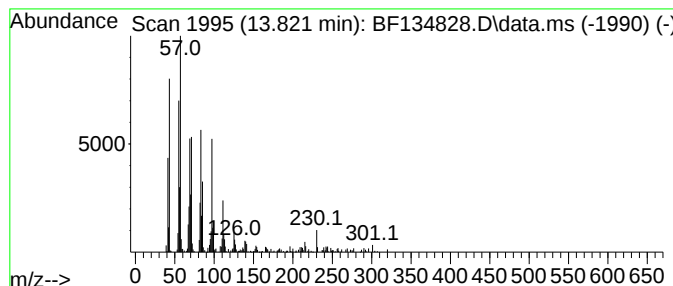
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecafluorooctanoic aci... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.821	3.33 ng	229772	Chrysene-d12	13.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	95
2			1-Pentadecanethiol	244	C15H32S	025276-70-4	91
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87
5			Cyclotetradecane	196	C14H28	000295-17-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081723\
Data File : BF134828.D
Acq On : 17 Aug 2023 21:05
Operator : CG\JU
Sample : 04042-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

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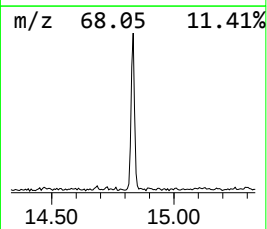
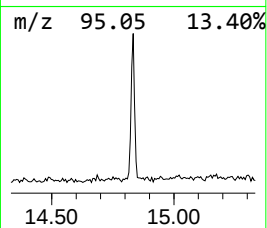
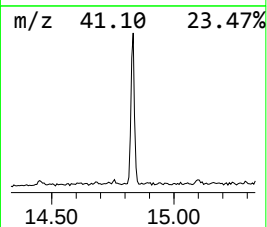
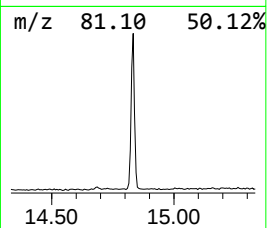
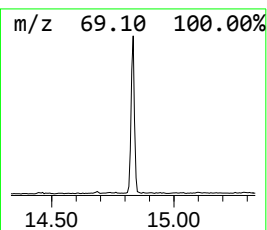
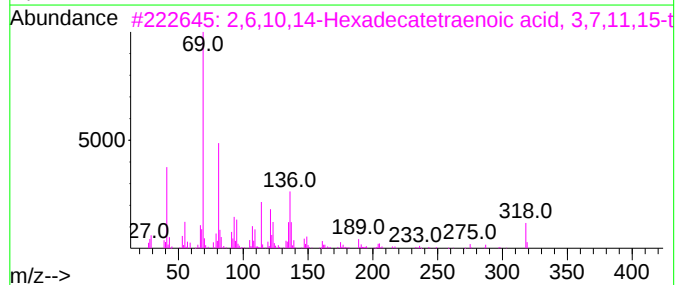
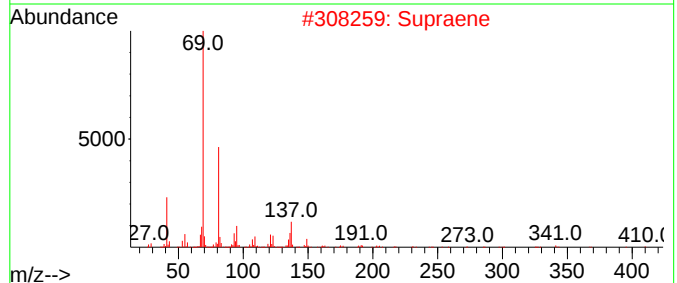
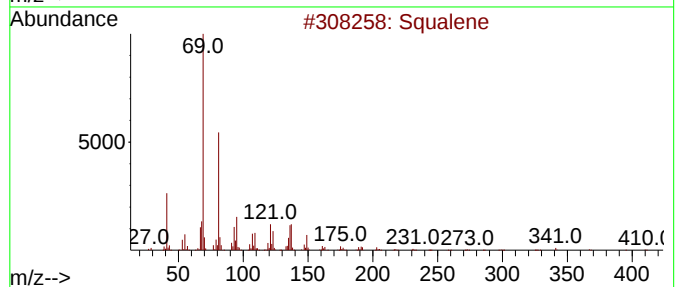
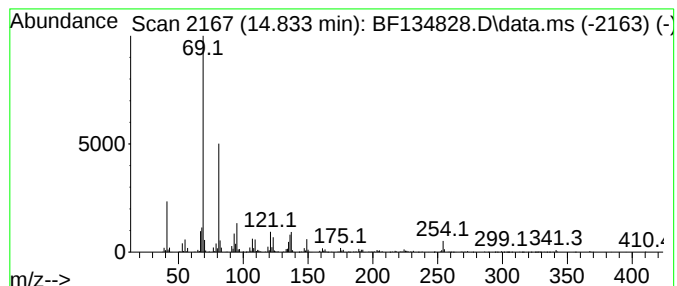
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Squalene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.833	12.00 ng	641528	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	95
2			Supraene	410	C30H50	007683-64-9	80
3			2,6,10,14-Hexadecatetraenoic aci...	318	C21H34O2	042207-88-5	72
4			2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	59
5			Trifluoroacetyl-lavandulol	250	C12H17F3O2	028673-24-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081723\
Data File : BF134828.D
Acq On : 17 Aug 2023 21:05
Operator : CG\JU
Sample : 04042-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-376

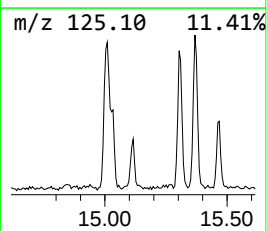
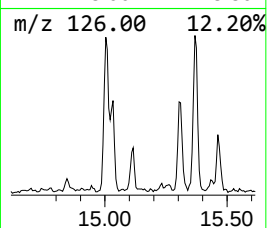
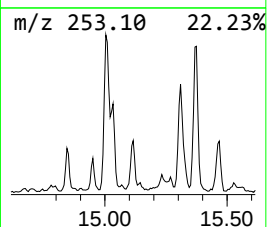
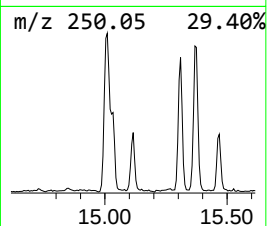
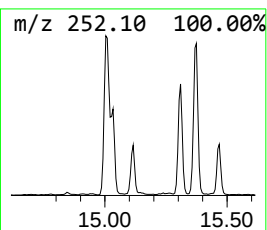
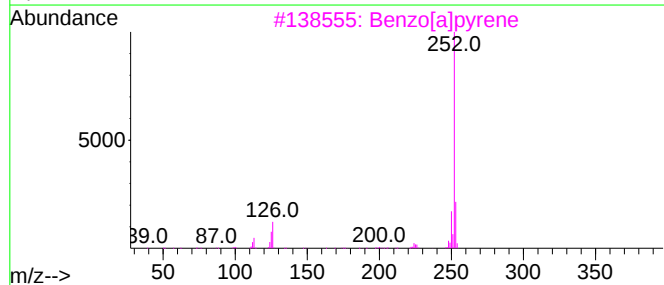
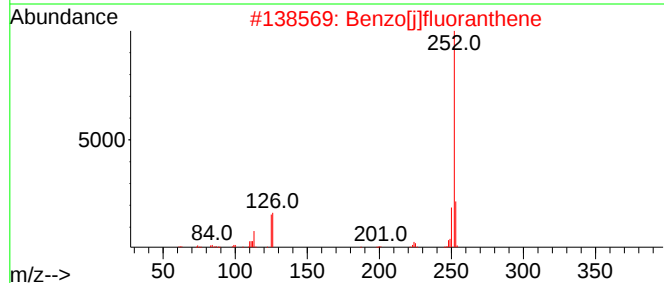
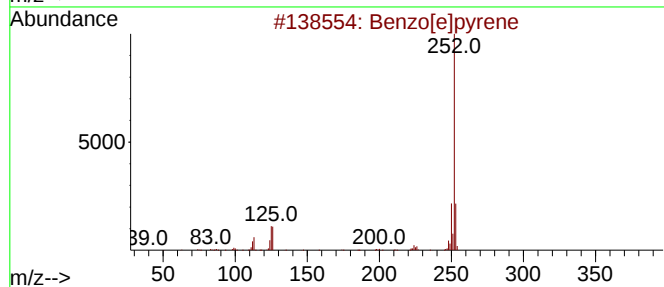
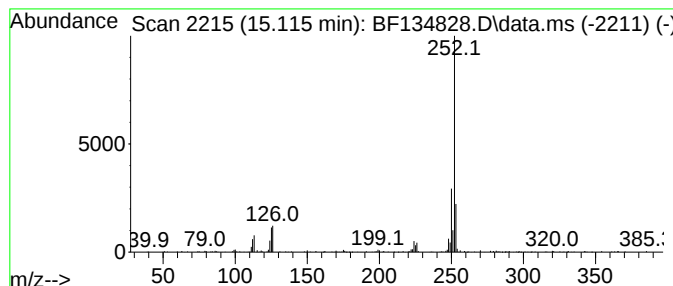
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.115	3.07 ng	164308	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	96
2			Benzo[j]fluoranthene	252	C20H12	000205-82-3	93
3			Benzo[a]pyrene	252	C20H12	000050-32-8	93
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5			Perylene	252	C20H12	000198-55-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081723\
Data File : BF134828.D
Acq On : 17 Aug 2023 21:05
Operator : CG\JU
Sample : 04042-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-376

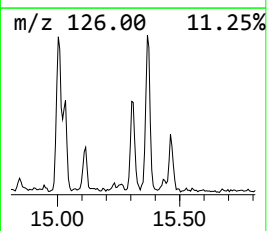
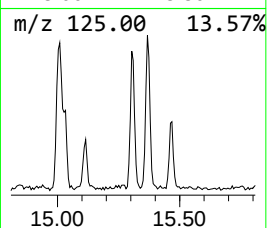
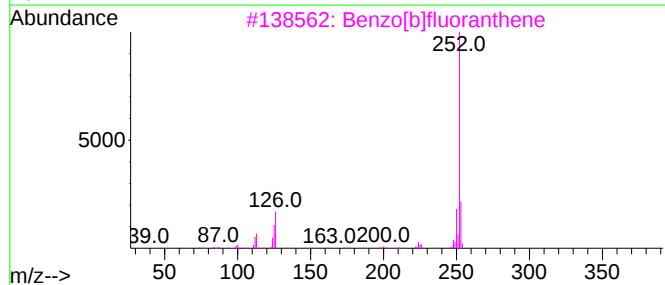
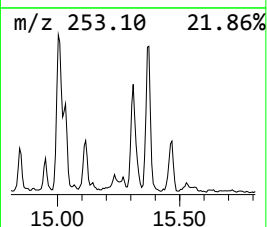
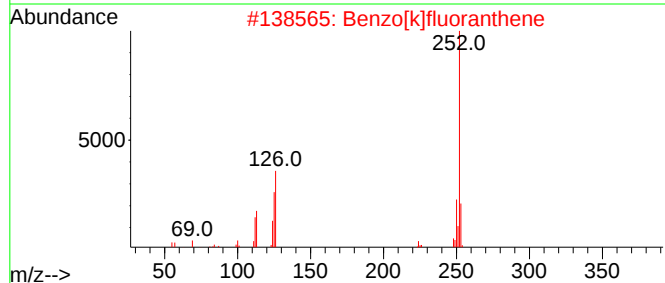
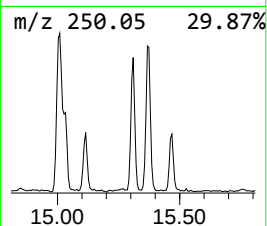
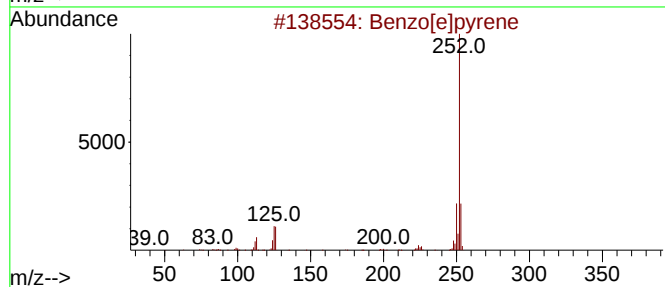
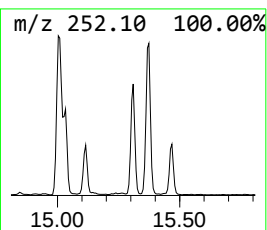
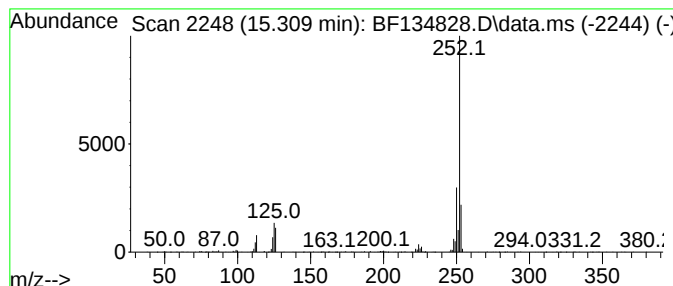
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[j]fluoranthene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.309	7.10 ng	379474	Perylene-d12	15.439

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	95
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90
5			Benzo[a]pyrene	252	C20H12	000050-32-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081723\
Data File : BF134828.D
Acq On : 17 Aug 2023 21:05
Operator : CG\JU
Sample : 04042-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.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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8,9-Dihydrocycl...	11.927	3.3	ng	155079	4	11.316	931974	20.0
Phenanthrene, 2...	12.368	2.1	ng	98818	4	11.316	931974	20.0
Octadecanoic acid	12.639	2.8	ng	132008	4	11.316	931974	20.0
9-Octadecenamid...	13.398	10.6	ng	732284	5	13.963	1379390	20.0
Pentadecafluoro...	13.821	3.3	ng	229772	5	13.963	1379390	20.0
Squalene	14.833	12.0	ng	641528	6	15.439	1069010	20.0
Benzo[e]pyrene	15.115	3.1	ng	164308	6	15.439	1069010	20.0
Benzo[j]fluoran...	15.309	7.1	ng	379474	6	15.439	1069010	20.0