

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031822\
 Data File : BF127409.D
 Acq On : 18 Mar 2022 23:44
 Operator : CG\JU
 Sample : N1983-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 124042

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-BF031022.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127409.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.481	539	543	553	rBV	1661183	1514347	100.00%	13.040%
2	6.492	710	715	718	rBV	1403962	1335161	88.17%	11.497%
3	6.857	773	777	783	rBV	288589	253741	16.76%	2.185%
4	7.422	869	873	876	rBV	926948	911583	60.20%	7.849%
5	8.110	986	990	992	rBV	21823	19759	1.30%	0.170%
6	8.139	992	995	998	rBV	389979	311220	20.55%	2.680%
7	8.716	1090	1093	1098	rVB	48772	35902	2.37%	0.309%
8	9.216	1173	1178	1181	rBV	1840049	1452604	95.92%	12.508%
9	9.280	1185	1189	1192	rBV	56985	46341	3.06%	0.399%
10	9.475	1219	1222	1227	rBV3	16347	24111	1.59%	0.208%
11	9.551	1231	1235	1238	rVV2	22895	24288	1.60%	0.209%
12	9.669	1252	1255	1260	rVB3	16295	21520	1.42%	0.185%
13	9.810	1275	1279	1284	rVB	45915	42062	2.78%	0.362%
14	9.892	1290	1293	1297	rVV	428898	383085	25.30%	3.299%
15	9.927	1297	1299	1302	rVB	52394	37167	2.45%	0.320%
16	10.098	1325	1328	1331	rVB2	37431	36157	2.39%	0.311%
17	10.133	1331	1334	1338	rBV3	21955	21959	1.45%	0.189%
18	10.310	1358	1364	1367	rBV2	49859	55746	3.68%	0.480%
19	10.445	1382	1387	1393	rBV	46581	46022	3.04%	0.396%
20	10.527	1396	1401	1404	rVV2	24259	30670	2.03%	0.264%
21	10.692	1424	1429	1432	rBV	1278123	1138371	75.17%	9.802%
22	10.786	1442	1445	1450	rBV	54878	68092	4.50%	0.586%
23	10.992	1475	1480	1482	rBV5	13567	21742	1.44%	0.187%
24	11.233	1518	1521	1524	rBV	49616	40155	2.65%	0.346%
25	11.386	1543	1547	1549	rBV	525335	428867	28.32%	3.693%
26	11.410	1549	1551	1554	rVV	200643	150363	9.93%	1.295%
27	11.463	1554	1560	1564	rVB	27861	31739	2.10%	0.273%
28	11.621	1582	1587	1591	rBV2	19769	31352	2.07%	0.270%
29	11.663	1591	1594	1596	rVB	36915	30862	2.04%	0.266%
30	11.763	1608	1611	1614	rBV	26574	23325	1.54%	0.201%
31	11.886	1627	1632	1634	rBV	34387	38701	2.56%	0.333%
32	11.916	1634	1637	1641	rVB2	57383	76104	5.03%	0.655%
33	11.998	1648	1651	1654	rBV2	40511	44786	2.96%	0.386%
34	12.068	1660	1663	1666	rBV	40925	34213	2.26%	0.295%
35	12.327	1702	1707	1710	rBV2	12351	19879	1.31%	0.171%
36	12.433	1722	1725	1727	rBV	30583	30019	1.98%	0.258%

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Stop Thrs : 0

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.457	1727	1729	1731	rBV	25159	25733	1.70%	0.222%
38	12.604	1751	1754	1758	rBV	153881	135716	8.96%	1.169%
39	12.833	1790	1793	1795	rVV	111367	78733	5.20%	0.678%
40	12.886	1799	1802	1809	rVB	32695	39977	2.64%	0.344%
41	12.974	1812	1817	1820	rBV	1756279	1442955	95.29%	12.425%
42	13.186	1851	1853	1855	rVB	33168	24076	1.59%	0.207%
43	13.268	1862	1867	1869	rBV3	18972	23380	1.54%	0.201%
44	13.527	1908	1911	1914	rBV2	29622	25956	1.71%	0.223%
45	13.857	1962	1967	1970	rBV4	29592	33287	2.20%	0.287%
46	13.998	1987	1991	1993	rVB	71500	61414	4.06%	0.529%
47	14.033	1993	1997	2000	rBV	358230	321657	21.24%	2.770%
48	14.115	2008	2011	2017	rVB6	17257	23758	1.57%	0.205%
49	14.174	2017	2021	2023	rBV	28195	28336	1.87%	0.244%
50	14.421	2061	2063	2066	rVB	28947	22049	1.46%	0.190%
51	14.480	2071	2073	2078	rVB2	23944	24451	1.61%	0.211%
52	14.780	2121	2124	2127	rVB3	18505	22153	1.46%	0.191%
53	15.127	2179	2183	2186	rVB2	19980	25800	1.70%	0.222%
54	15.456	2236	2239	2244	rVB	16817	21948	1.45%	0.189%
55	15.515	2244	2249	2256	rBV	241854	322076	21.27%	2.773%
56	15.945	2318	2322	2330	rVB3	21254	38139	2.52%	0.328%
57	16.451	2404	2408	2414	rVB2	14747	24475	1.62%	0.211%
58	17.756	2620	2630	2636	rBV9	11173	35377	2.34%	0.305%

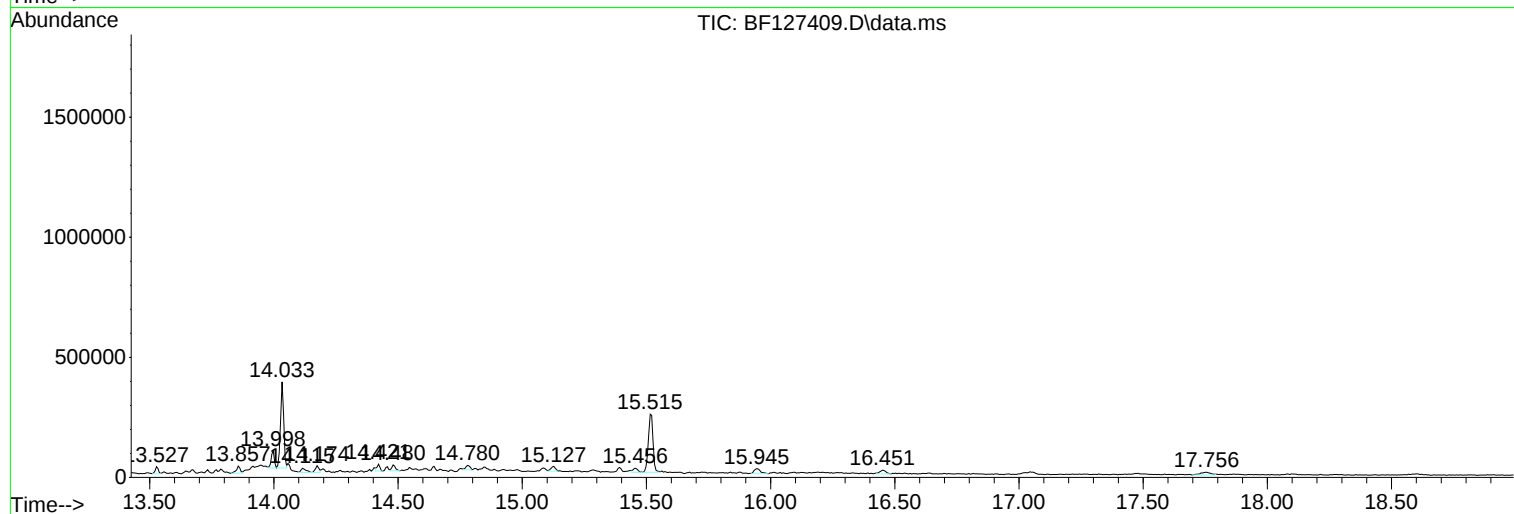
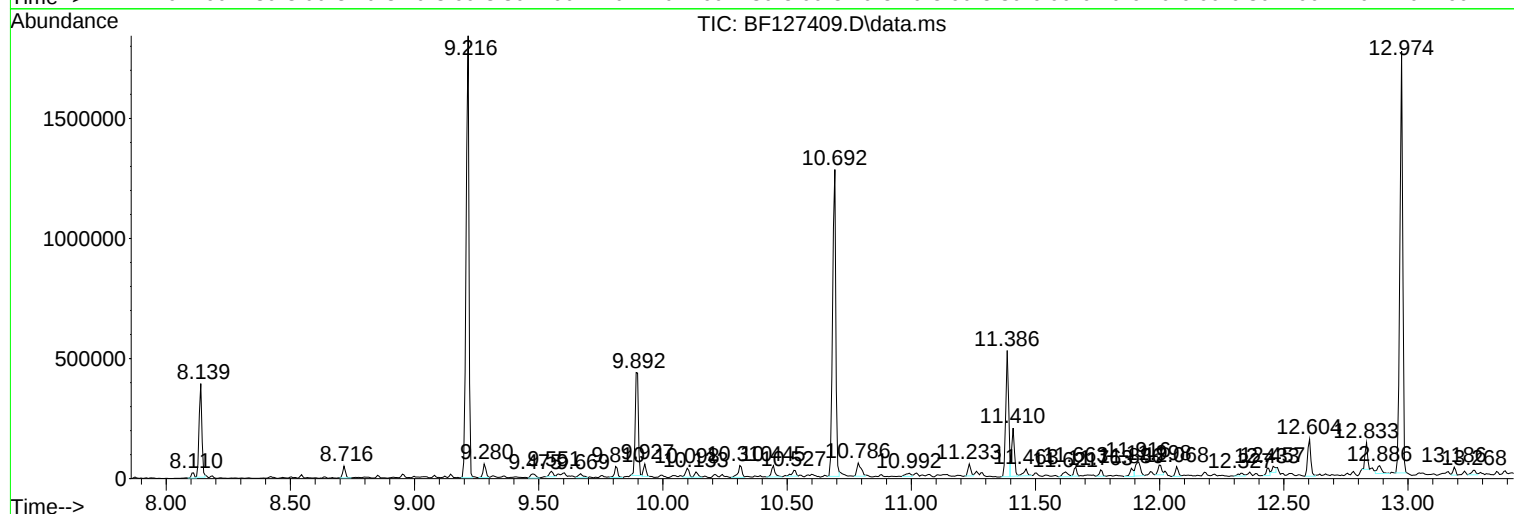
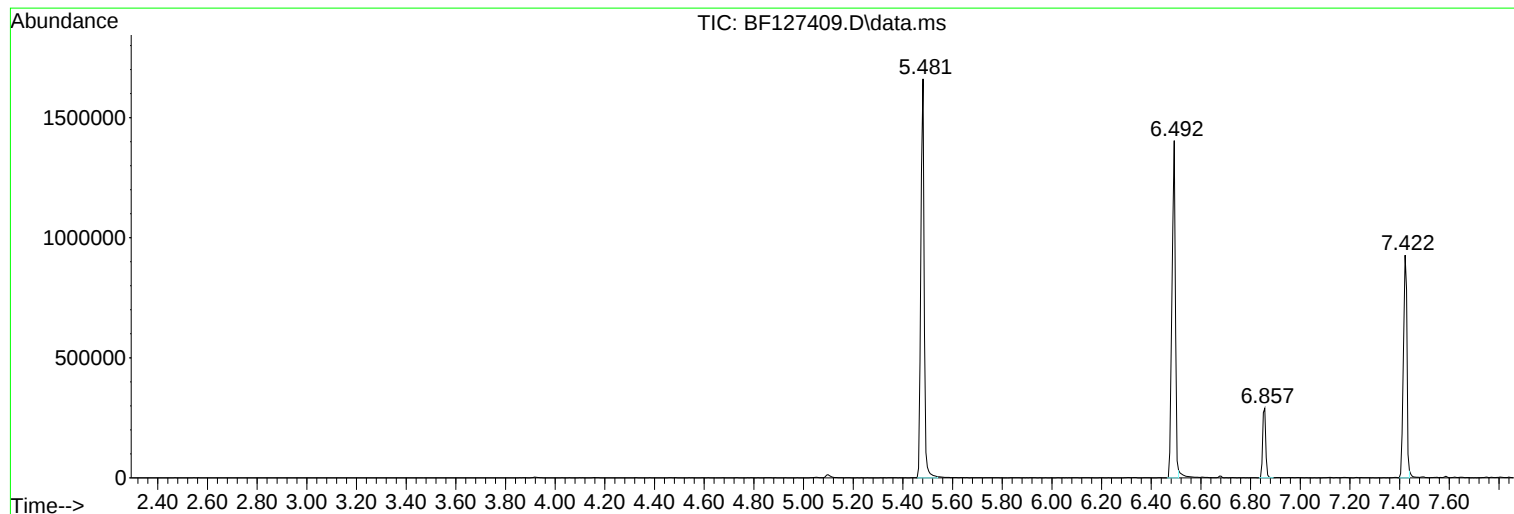
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Operator : CG\JU
Sample : N1983-03
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
124042

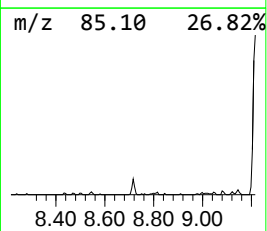
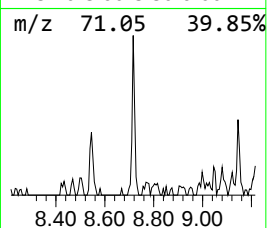
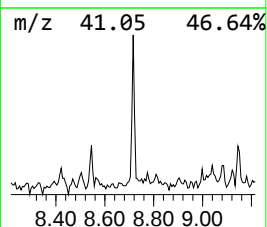
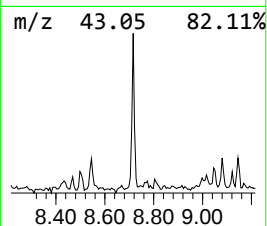
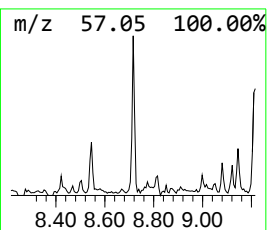
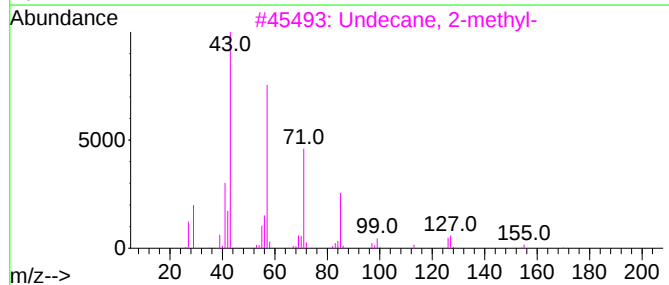
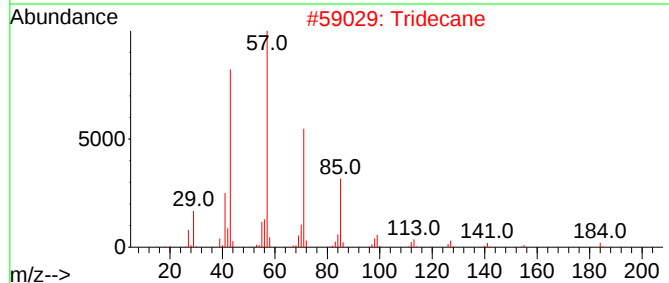
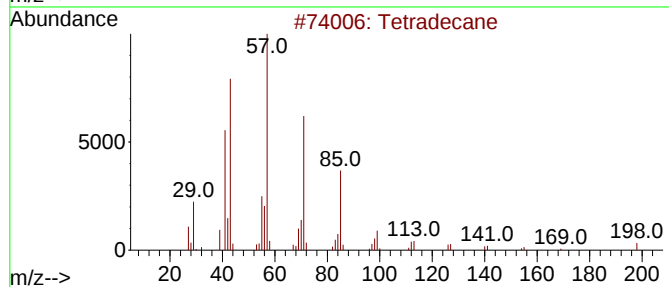
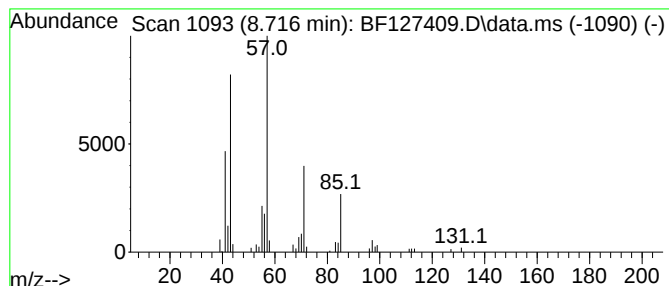
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Tetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.716	2.31 ng	35902	Naphthalene-d8	8.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	80
2			Tridecane	184	C13H28	000629-50-5	72
3			Undecane, 2-methyl-	170	C12H26	007045-71-8	72
4			Hexadecane	226	C16H34	000544-76-3	64
5			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	64



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124042

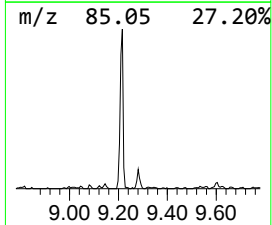
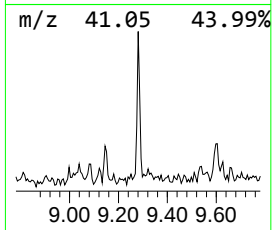
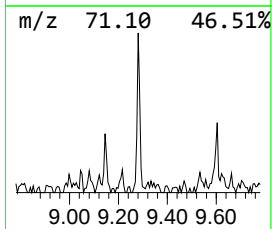
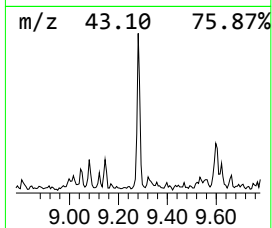
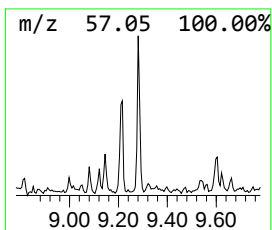
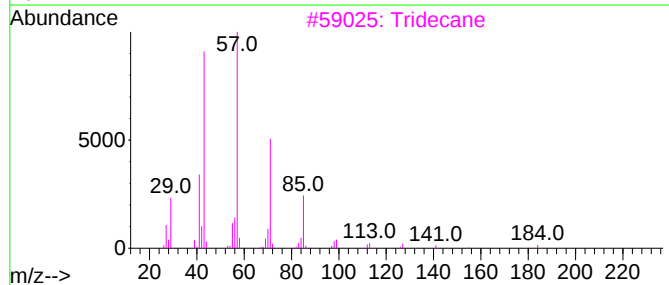
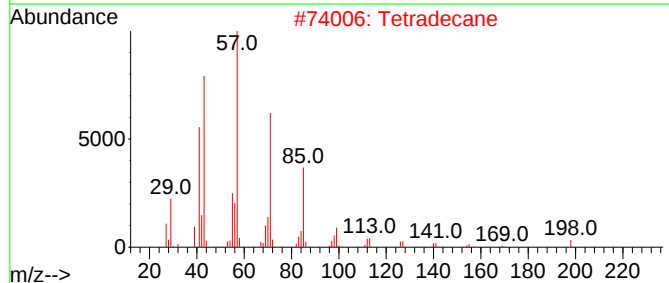
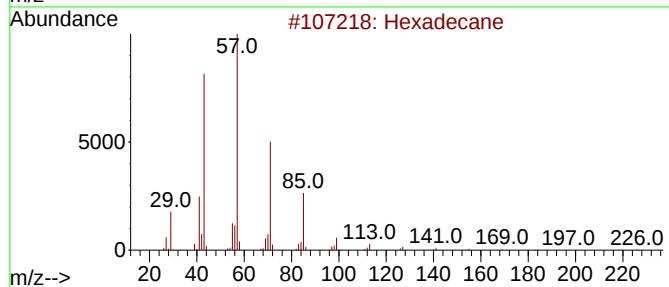
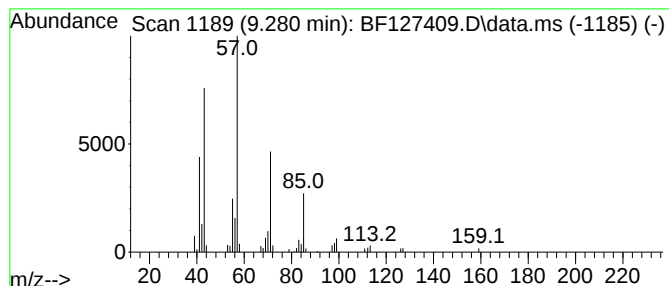
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.280	2.42 ng	46341	Acenaphthene-d10	9.898

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	91
2		Tetradecane	198	C14H30	000629-59-4	86
3		Tridecane	184	C13H28	000629-50-5	86
4		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	86
5		Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	83



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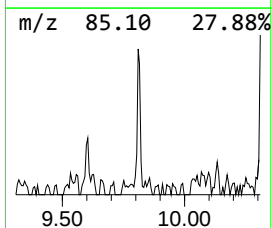
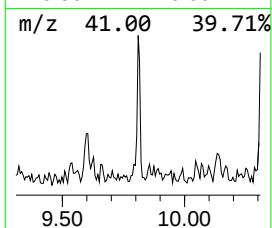
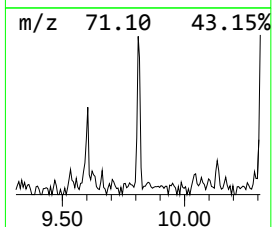
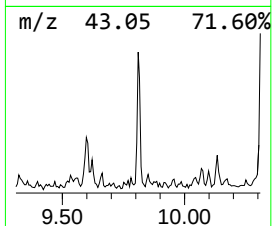
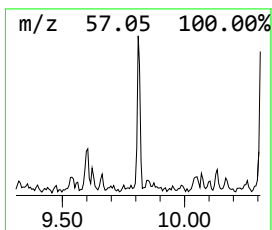
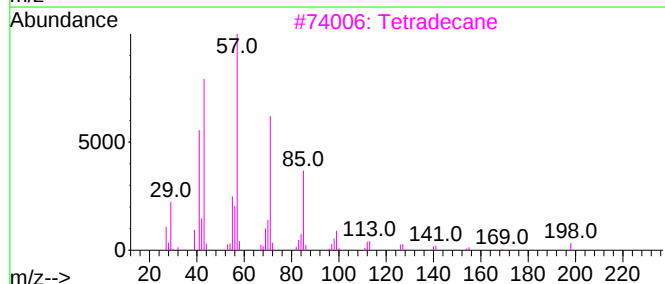
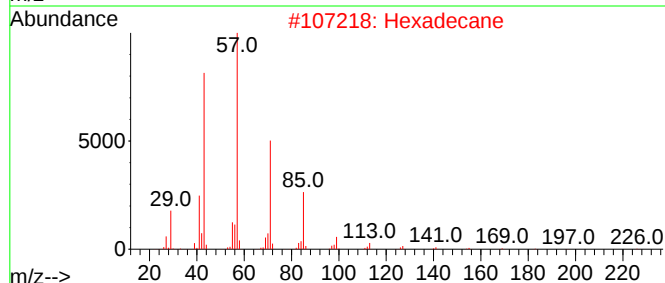
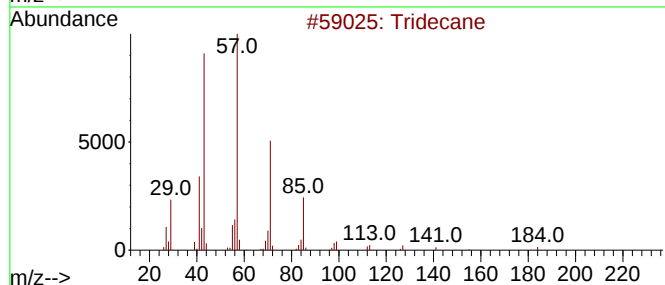
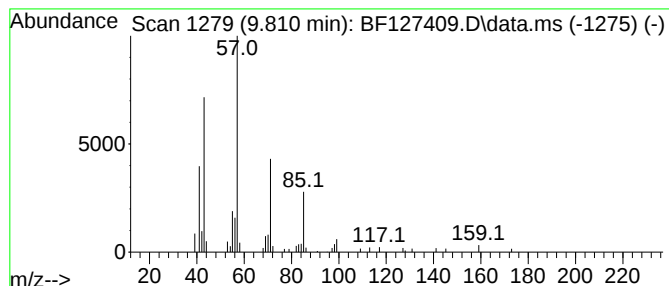
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Tridecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.810	2.20 ng	42062	Acenaphthene-d10	9.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	90
2			Hexadecane	226	C16H34	000544-76-3	87
3			Tetradecane	198	C14H30	000629-59-4	72
4			Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	72
5			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64



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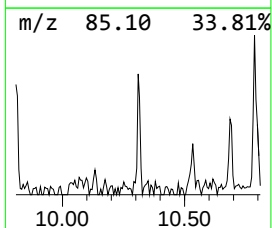
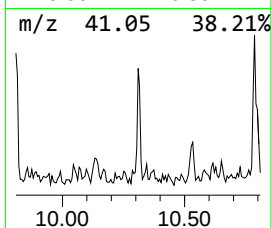
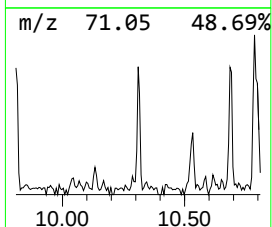
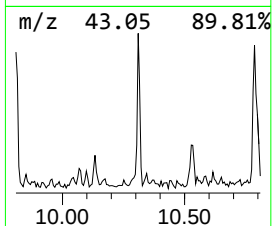
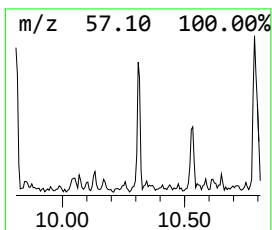
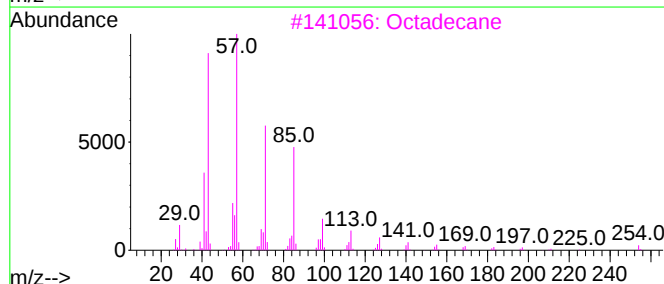
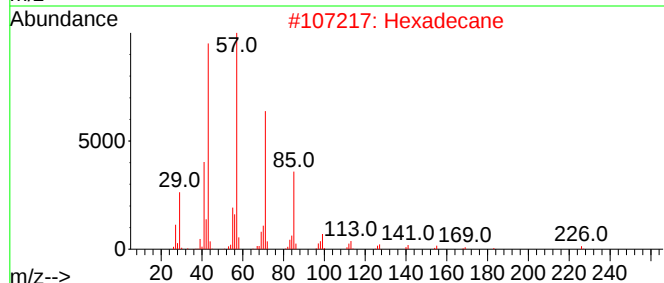
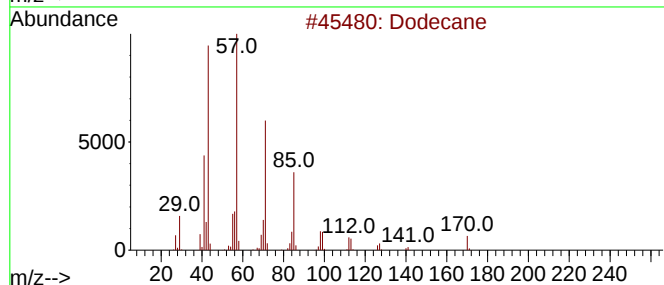
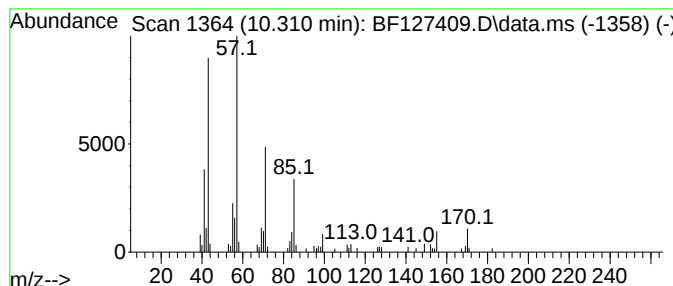
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.310	2.91 ng	55746	Acenaphthene-d10	9.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	81
2			Hexadecane	226	C16H34	000544-76-3	68
3			Octadecane	254	C18H38	000593-45-3	58
4			Tridecane	184	C13H28	000629-50-5	58
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	58



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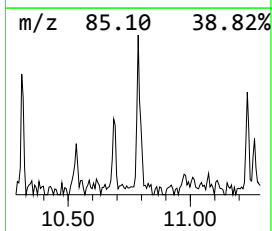
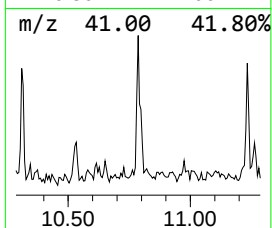
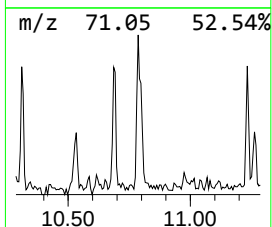
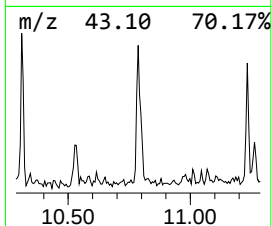
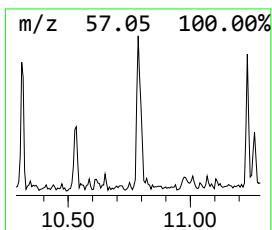
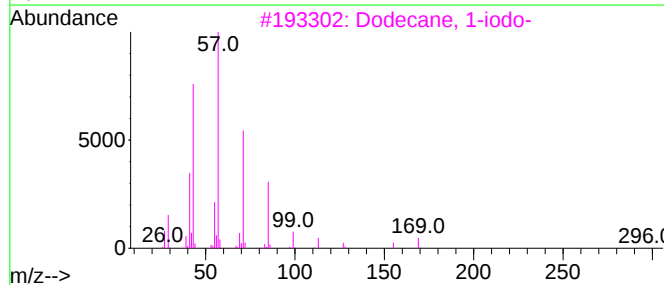
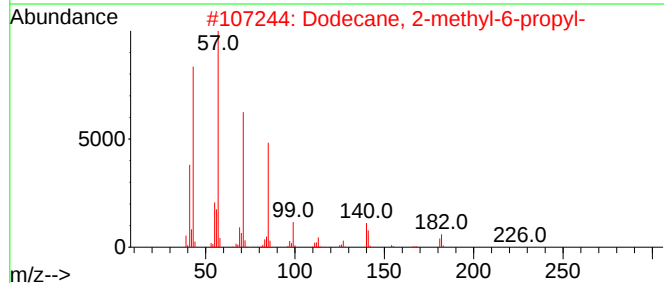
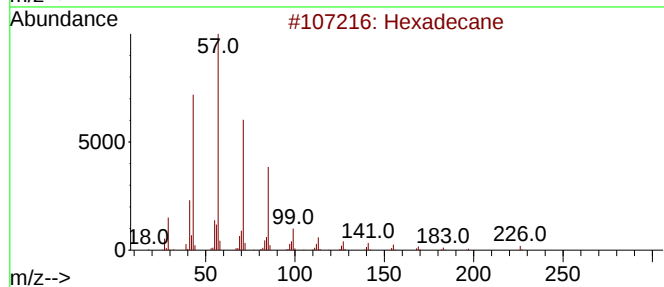
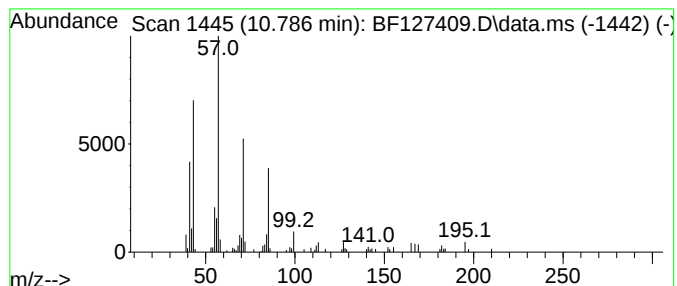
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Dodecane, 2-methyl-6-propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.786	3.18 ng	68092	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	72
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	72
3			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
4			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	72
5			Heptadecane	240	C17H36	000629-78-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031822\
Data File : BF127409.D
Acq On : 18 Mar 2022 23:44
Operator : CG\JU
Sample : N1983-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

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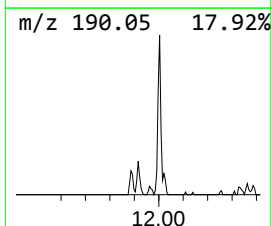
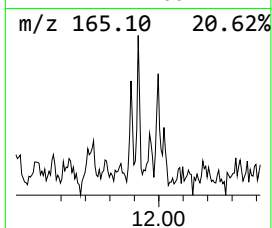
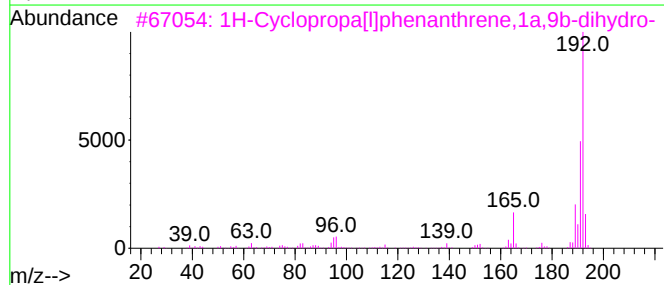
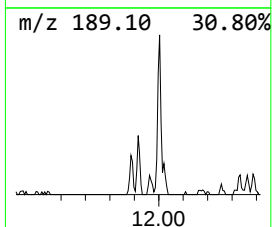
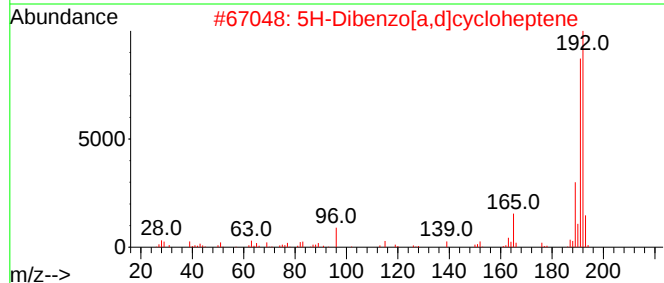
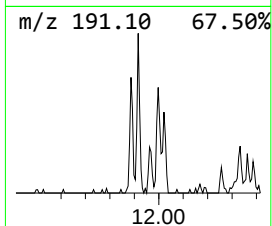
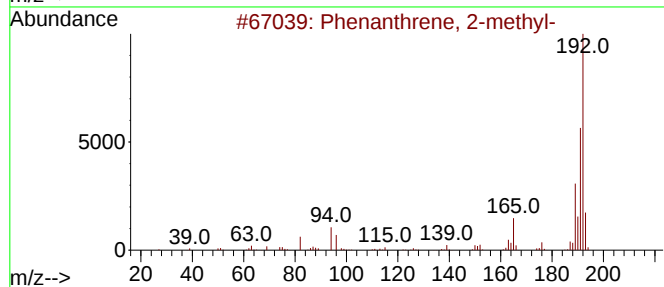
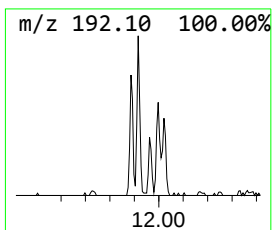
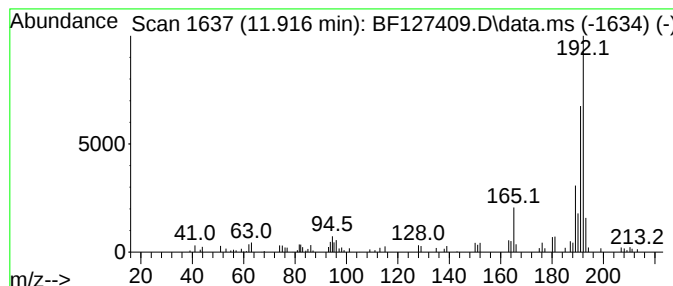
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	3.55 ng	76104	Phenanthrene-d10	11.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
2		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	94
3		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031822\
Data File : BF127409.D
Acq On : 18 Mar 2022 23:44
Operator : CG\JU
Sample : N1983-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

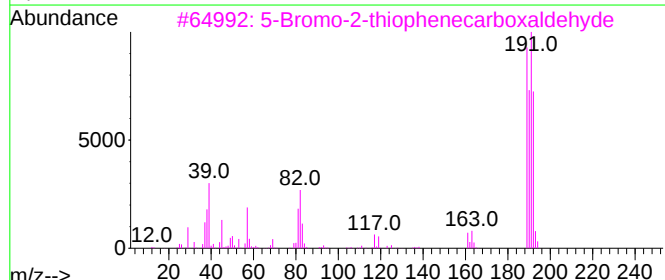
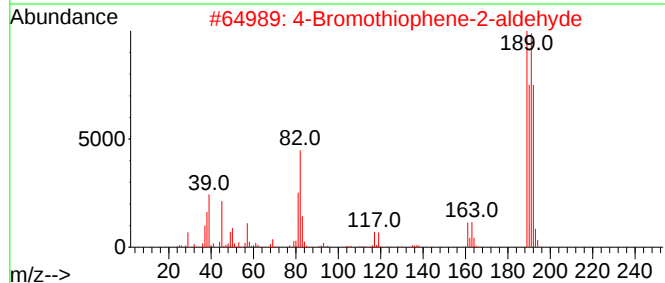
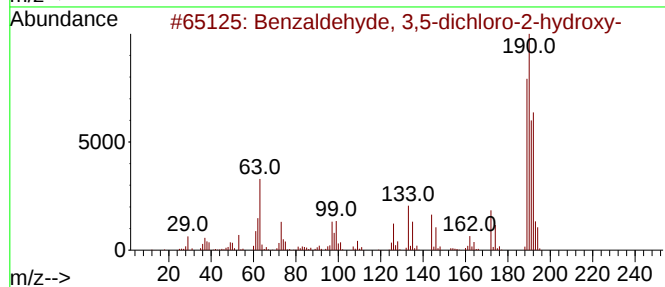
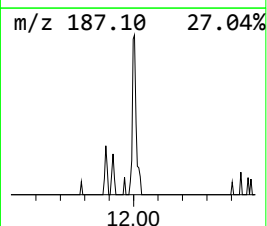
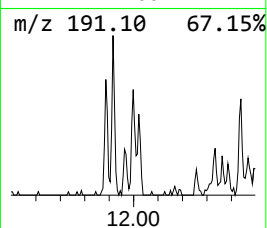
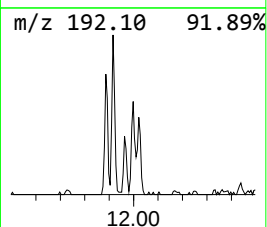
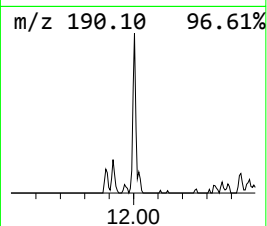
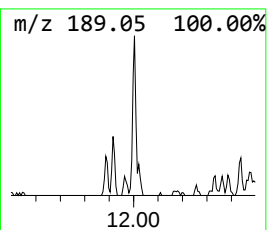
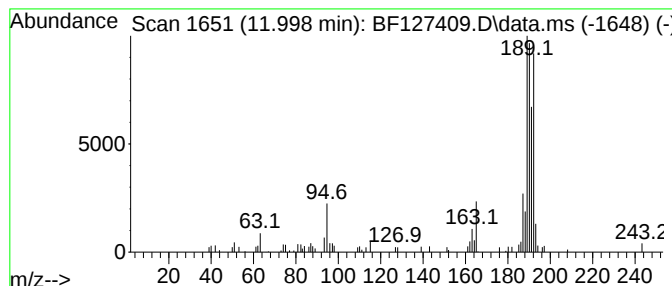
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzaldehyde, 3,5-dichloro-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.998	2.09 ng	44786	Phenanthrene-d10		11.386	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...		190	C7H4Cl2O2	000090-60-8	72
2	4-Bromothiophene-2-aldehyde		190	C5H3BrOS	018791-75-8	58
3	5-Bromo-2-thiophenecarboxaldehyde		190	C5H3BrOS	004701-17-1	52
4	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	47
5	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031822\
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Sample : N1983-03
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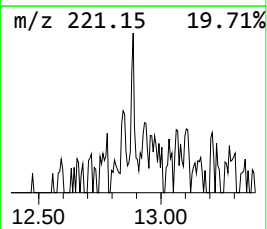
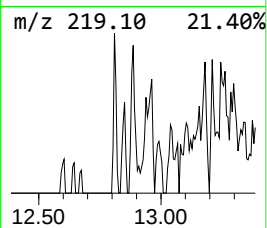
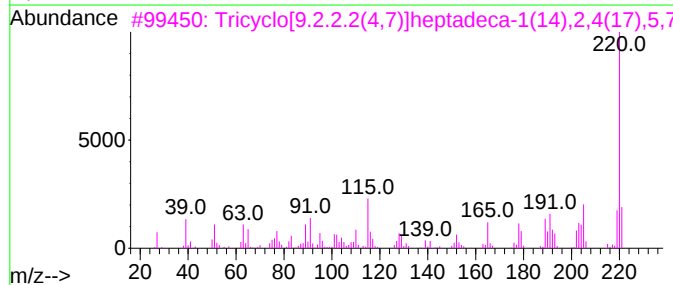
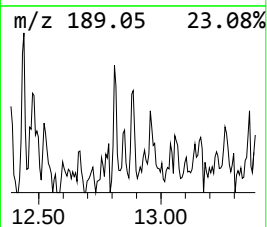
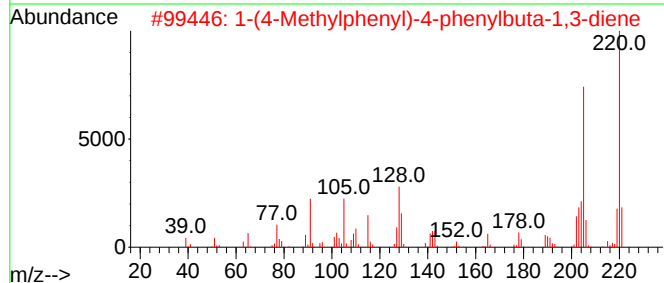
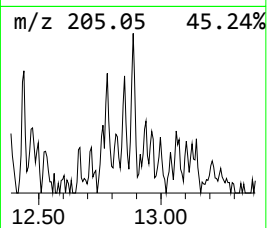
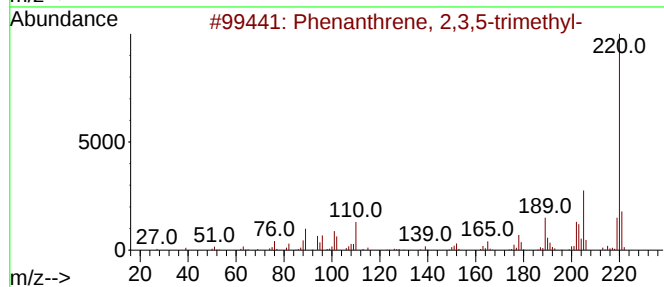
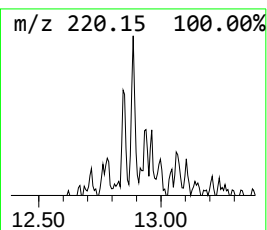
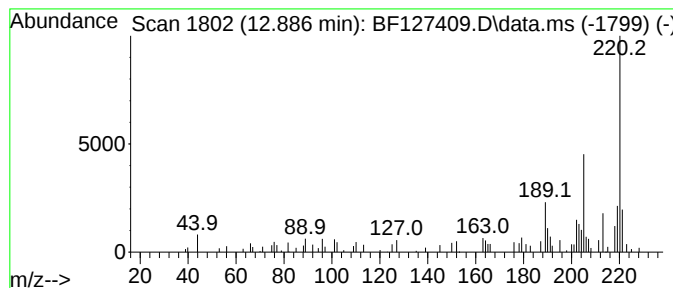
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.886	2.49 ng	39977	Chrysene-d12	14.033	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	90
2	1-(4-Methylphenyl)-4-phenylbuta-...	220	C17H16	037985-11-8	59
3	Tricyclo[9.2.2.2(4,7)]heptadeca-...	220	C17H16	049576-90-1	53
4	4-Phenyl-2-naphthol	220	C16H12O	036159-74-7	53
5	Benzo[b]dihdropyran, 6-hydroxy-...	220	C14H20O2	050442-70-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031822\
Data File : BF127409.D
Acq On : 18 Mar 2022 23:44
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Sample : N1983-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

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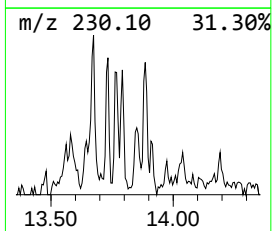
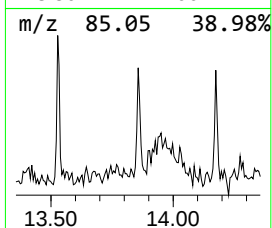
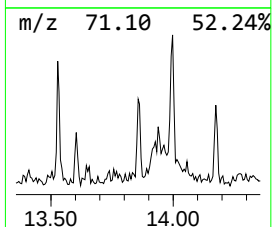
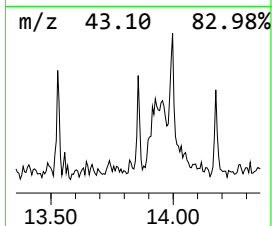
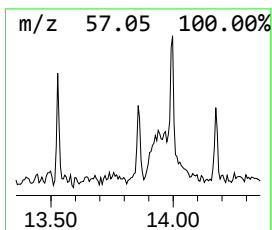
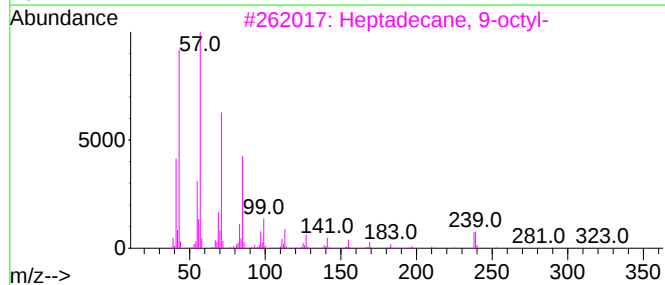
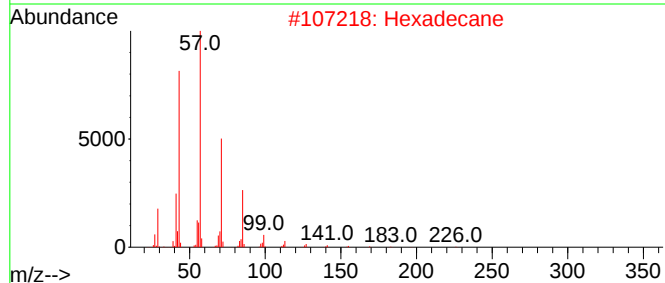
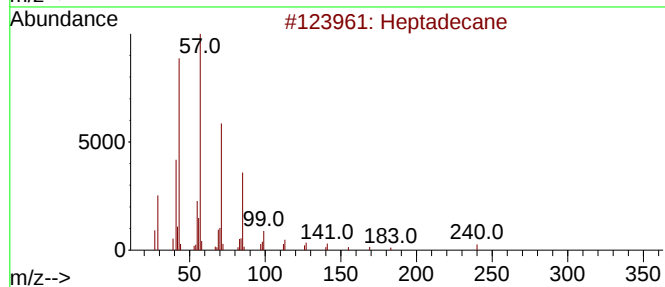
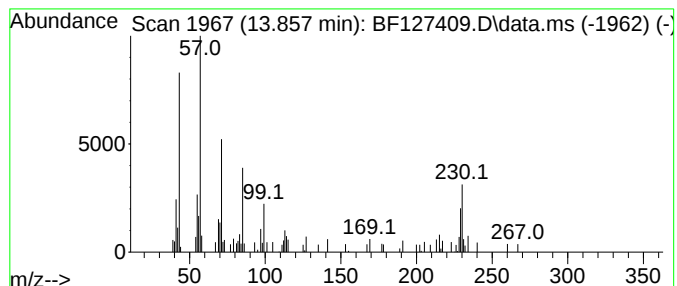
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	2.07 ng	33287	Chrysene-d12	14.033

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	55
2		Hexadecane	226	C16H34	000544-76-3	55
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	49
4		Hexacosane	366	C26H54	000630-01-3	49
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031822\
Data File : BF127409.D
Acq On : 18 Mar 2022 23:44
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Sample : N1983-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

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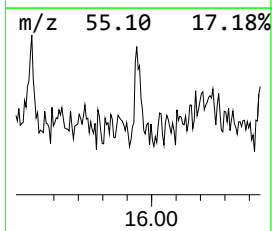
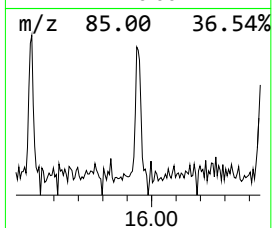
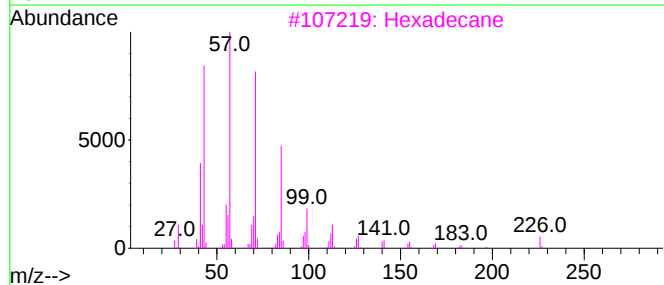
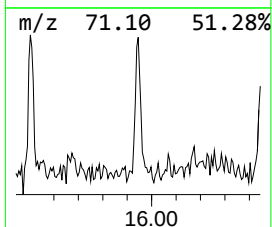
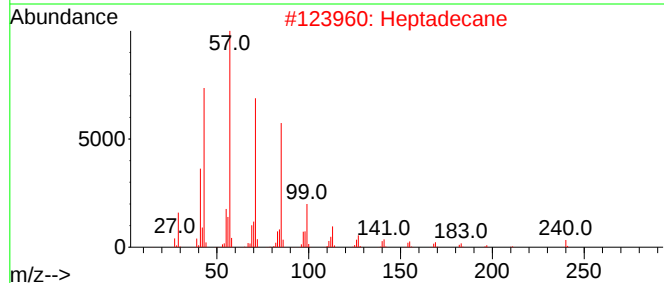
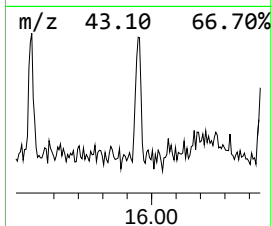
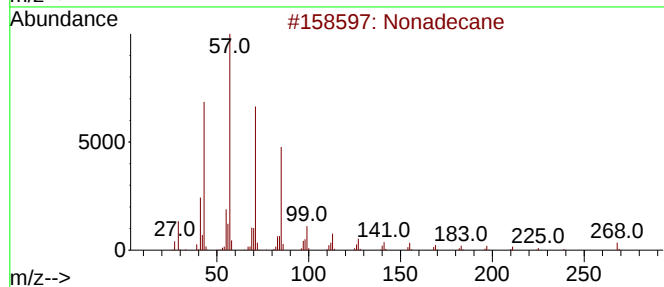
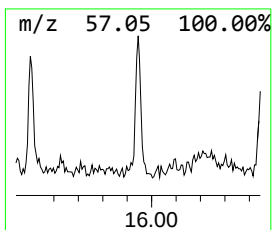
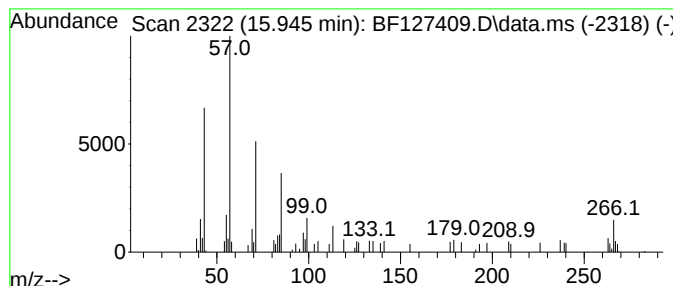
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	2.37 ng	38139	Perylene-d12	15.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	64
2			Heptadecane	240	C17H36	000629-78-7	64
3			Hexadecane	226	C16H34	000544-76-3	60
4			Pentacosane	352	C25H52	000629-99-2	49
5			Hexacosane	366	C26H54	000630-01-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031822\
Data File : BF127409.D
Acq On : 18 Mar 2022 23:44
Operator : CG\JU
Sample : N1983-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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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Hexadecane	9.280	2.4	ng	46341	3	9.898	383085	20.0
Tridecane	9.810	2.2	ng	42062	3	9.898	383085	20.0
Dodecane	10.310	2.9	ng	55746	3	9.898	383085	20.0
Dodecane, 2-met...	10.786	3.2	ng	68092	4	11.386	428867	20.0
Phenanthrene, 2...	11.916	3.5	ng	76104	4	11.386	428867	20.0
Benzaldehyde, 3...	11.998	2.1	ng	44786	4	11.386	428867	20.0
Phenanthrene, 2...	12.886	2.5	ng	39977	5	14.033	321657	20.0
Heptadecane	13.857	2.1	ng	33287	5	14.033	321657	20.0
Nonadecane	15.945	2.4	ng	38139	6	15.521	322076	20.0
unknown17.756	17.756	2.2	ng	35377	6	15.521	322076	20.0