

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
 Data File : BF130601.D
 Acq On : 10 Oct 2022 14:10
 Operator : CG\JU
 Sample : N5046-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-2

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-BF091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130601.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.781	454	458	471	rVB	448664	599456	24.10%	3.038%
2	5.210	527	531	546	rBV	2231162	2152700	86.53%	10.911%
3	5.610	596	599	604	rVB	49045	43695	1.76%	0.221%
4	6.251	704	708	711	rBV	2097622	2128808	85.57%	10.790%
5	6.604	765	768	772	rBV	762294	589723	23.70%	2.989%
6	7.175	860	865	868	rBV	1665956	1411597	56.74%	7.155%
7	7.839	971	978	982	rBV4	30171	78240	3.14%	0.397%
8	7.886	982	986	989	rVV	997808	767280	30.84%	3.889%
9	8.698	1119	1124	1128	rBV2	34121	35175	1.41%	0.178%
10	8.963	1165	1169	1172	rBV	3046267	2487761	100.00%	12.610%
11	9.045	1178	1183	1187	rVB	51991	48599	1.95%	0.246%
12	9.092	1187	1191	1196	rBV	64977	78187	3.14%	0.396%
13	9.298	1222	1226	1228	rBV	39821	37574	1.51%	0.190%
14	9.498	1256	1260	1263	rBV	47608	48022	1.93%	0.243%
15	9.574	1270	1273	1276	rBV	67881	51218	2.06%	0.260%
16	9.633	1276	1283	1287	rBV	1061033	845616	33.99%	4.286%
17	10.074	1354	1358	1363	rVB2	73319	68901	2.77%	0.349%
18	10.180	1373	1376	1382	rVB4	35199	44980	1.81%	0.228%
19	10.292	1392	1395	1400	rVB	35595	32506	1.31%	0.165%
20	10.421	1413	1417	1429	rVB	1335908	1123754	45.17%	5.696%
21	10.545	1435	1438	1443	rBV	110211	139371	5.60%	0.706%
22	10.727	1465	1469	1471	rBV2	25632	34712	1.40%	0.176%
23	10.992	1511	1514	1516	rBV	71729	60207	2.42%	0.305%
24	11.016	1516	1518	1525	rVB2	50007	58492	2.35%	0.296%
25	11.116	1531	1535	1537	rBV	821383	697546	28.04%	3.536%
26	11.139	1537	1539	1541	rVV	267391	206738	8.31%	1.048%
27	11.186	1545	1547	1551	rVB	47425	41394	1.66%	0.210%
28	11.416	1583	1586	1589	rBV	66210	56551	2.27%	0.287%
29	11.521	1600	1604	1610	rVB3	44869	54946	2.21%	0.279%
30	11.616	1617	1620	1622	rBV	113219	88795	3.57%	0.450%
31	11.645	1622	1625	1631	rVV2	145808	229240	9.21%	1.162%
32	11.721	1635	1638	1641	rVV2	116818	134878	5.42%	0.684%
33	11.745	1641	1642	1648	rVB	84842	72250	2.90%	0.366%
34	11.821	1653	1655	1658	rVB	43981	34853	1.40%	0.177%
35	11.910	1667	1670	1672	rBV	35743	34200	1.37%	0.173%
36	11.939	1672	1675	1680	rVB2	38706	37717	1.52%	0.191%

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37	12.163	1709	1713	1716	rBV	91974	90416	3.63%	0.458%
38	12.198	1716	1719	1721	rVV2	59224	81125	3.26%	0.411%
39	12.221	1721	1723	1725	rVV2	114094	91458	3.68%	0.464%
40	12.251	1725	1728	1732	rVV3	36576	44260	1.78%	0.224%
41	12.321	1734	1740	1744	rVB	205939	197526	7.94%	1.001%
42	12.439	1754	1760	1773	rBV4	228086	405705	16.31%	2.056%
43	12.551	1774	1779	1781	rVV	238500	240655	9.67%	1.220%
44	12.580	1781	1784	1786	rVV4	35241	41369	1.66%	0.210%
45	12.610	1786	1789	1792	rVB2	32034	35327	1.42%	0.179%
46	12.698	1800	1804	1811	rVB	1760233	1499052	60.26%	7.598%
47	12.768	1813	1816	1820	rBV	35661	46744	1.88%	0.237%
48	12.880	1833	1835	1838	rVB	48433	39619	1.59%	0.201%
49	12.945	1839	1846	1849	rBV4	38801	57969	2.33%	0.294%
50	12.980	1849	1852	1855	rVB	59517	52623	2.12%	0.267%
51	13.074	1865	1868	1871	rBV	47127	40606	1.63%	0.206%
52	13.104	1871	1873	1876	rVV	45326	35302	1.42%	0.179%
53	13.280	1896	1903	1905	rBV6	25293	42227	1.70%	0.214%
54	13.386	1919	1921	1925	rVB	28719	29854	1.20%	0.151%
55	13.498	1934	1940	1942	rBV3	38120	53544	2.15%	0.271%
56	13.598	1955	1957	1960	rBV2	33093	32337	1.30%	0.164%
57	13.745	1977	1982	1985	rBV2	621067	666215	26.78%	3.377%
58	13.768	1985	1986	1990	rVB	113372	86889	3.49%	0.440%
59	14.133	2043	2048	2051	rVB	49707	50966	2.05%	0.258%
60	14.162	2051	2053	2056	rBV	32912	30226	1.21%	0.153%
61	14.227	2061	2064	2066	rVB2	39854	33265	1.34%	0.169%
62	14.257	2066	2069	2071	rBV2	35429	33123	1.33%	0.168%
63	14.521	2110	2114	2116	rBV2	20727	34360	1.38%	0.174%
64	14.745	2148	2152	2158	rBV2	94014	158952	6.39%	0.806%
65	15.015	2195	2198	2203	rVB	64015	65838	2.65%	0.334%
66	15.068	2204	2207	2212	rBV	99228	107888	4.34%	0.547%
67	15.127	2213	2217	2221	rBV	540972	579232	23.28%	2.936%
68	16.427	2434	2438	2445	rVB2	39595	68625	2.76%	0.348%

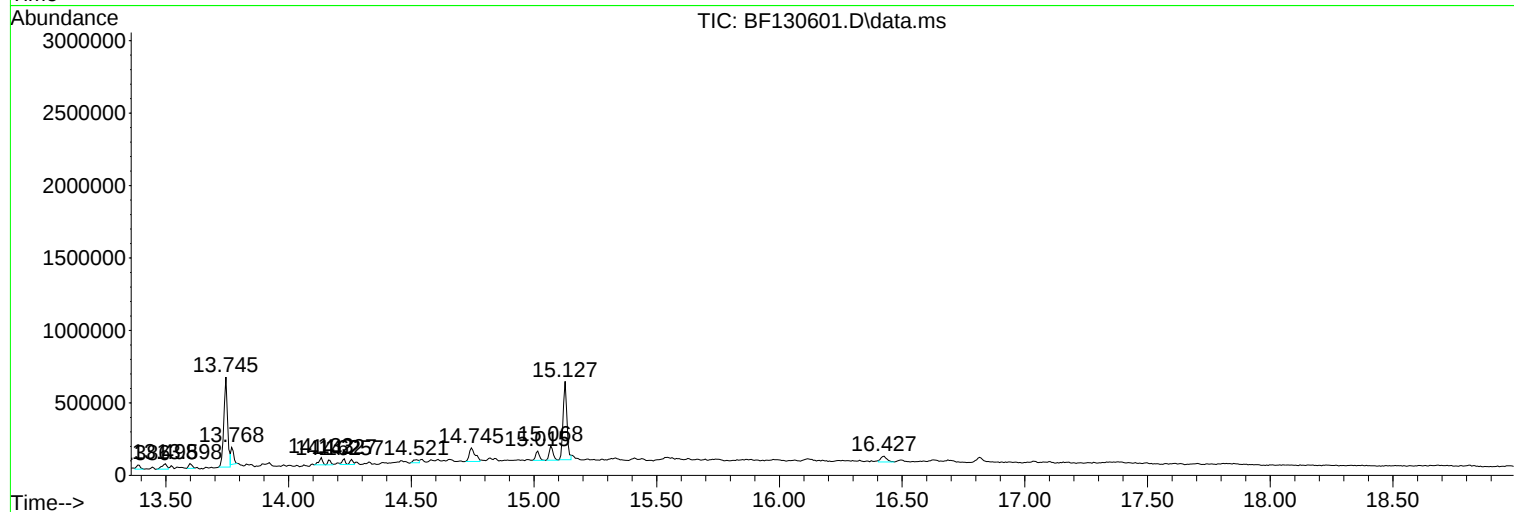
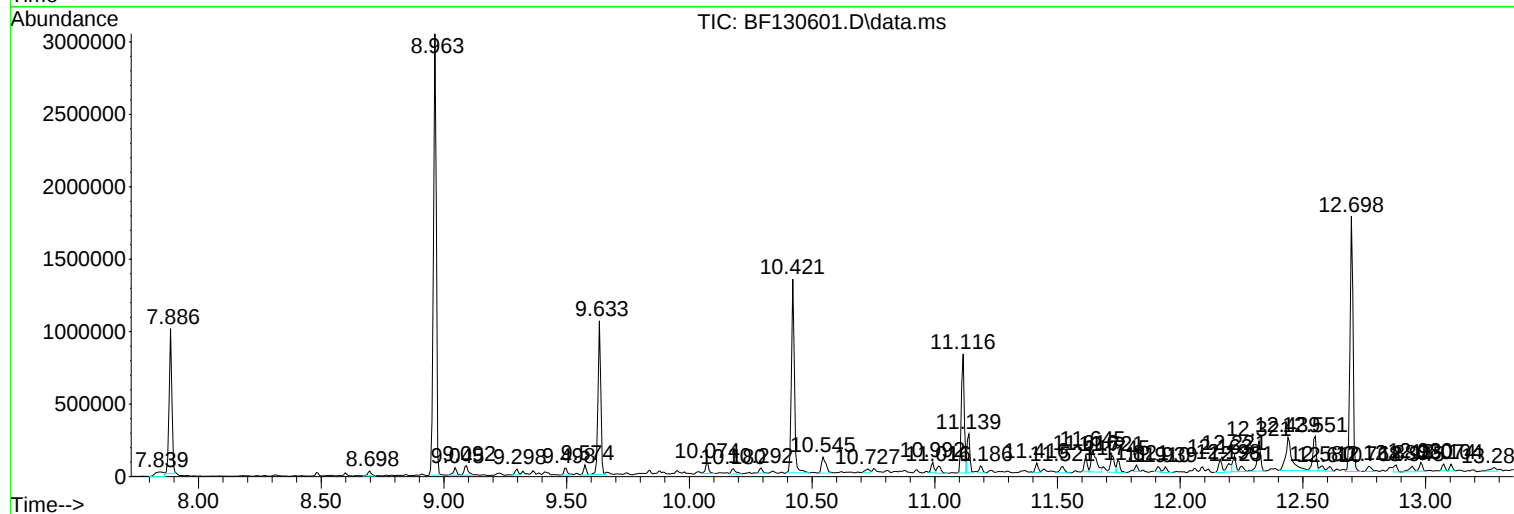
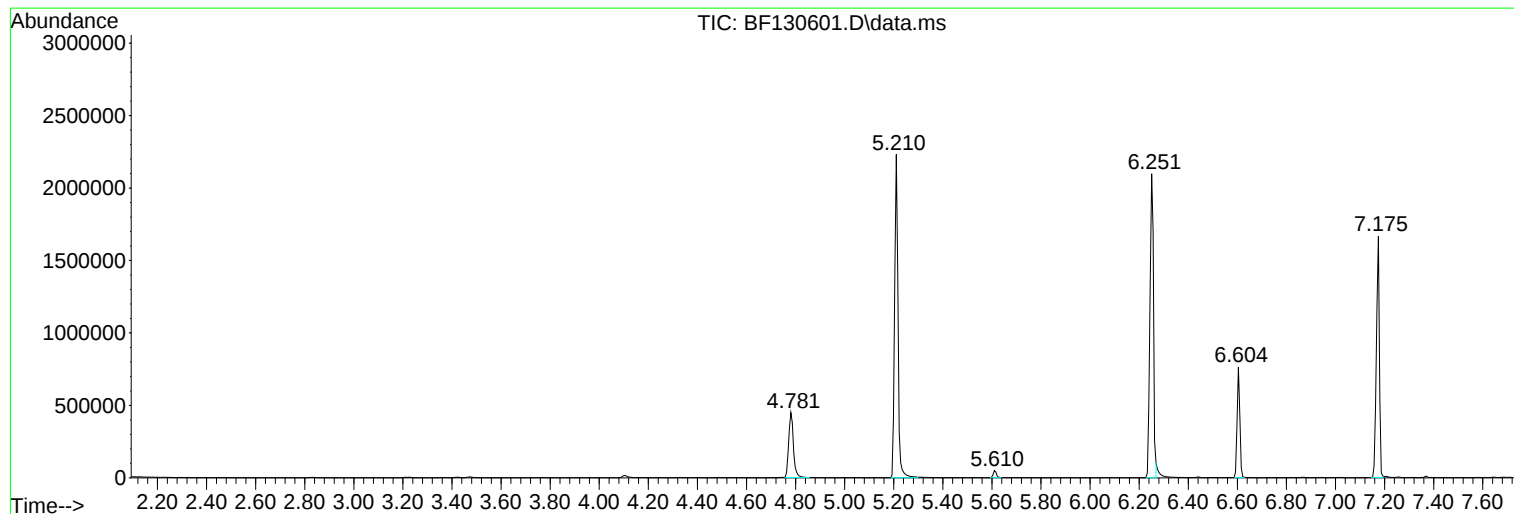
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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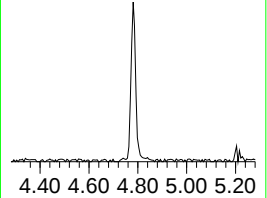
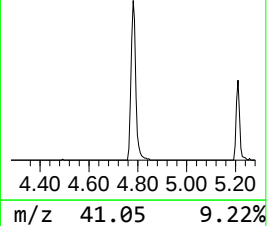
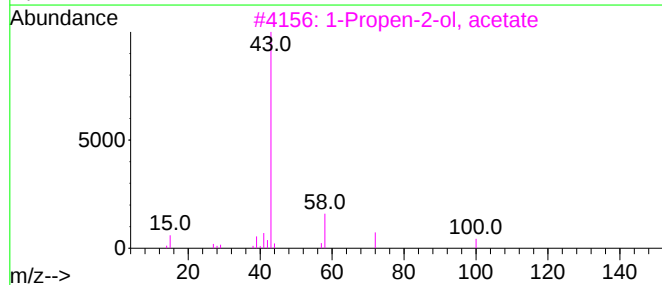
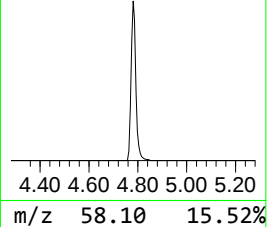
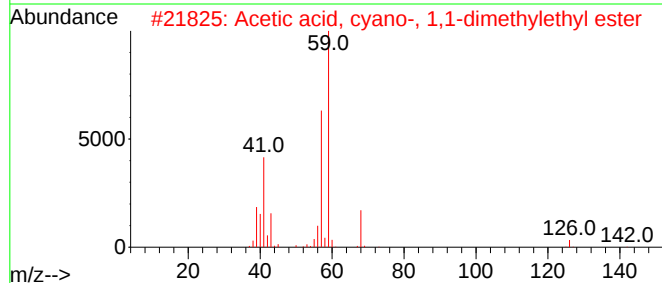
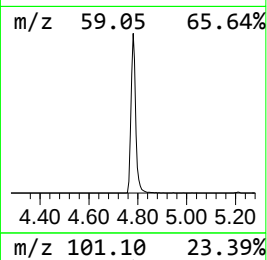
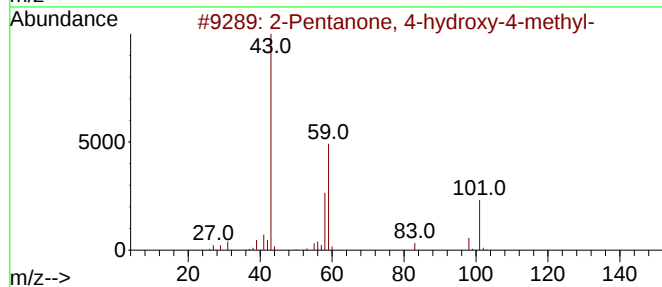
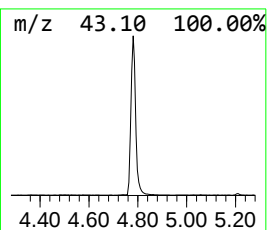
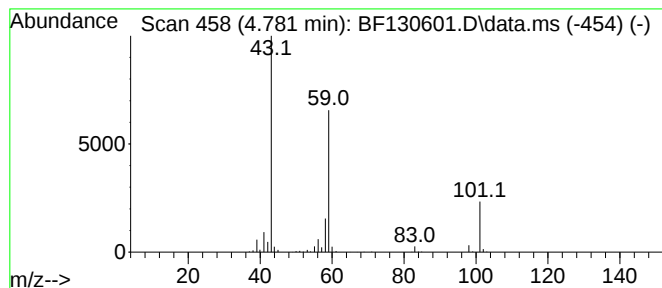
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.781	20.33 ng	599456	1,4-Dichlorobenzene-d4	6.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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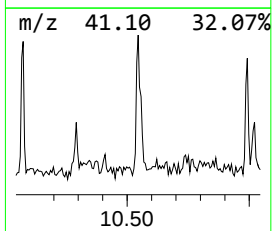
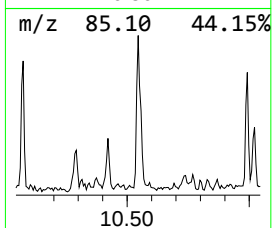
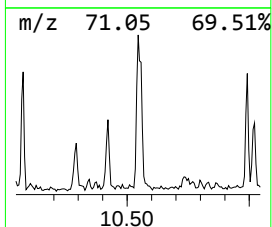
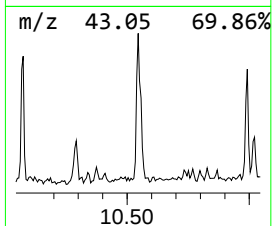
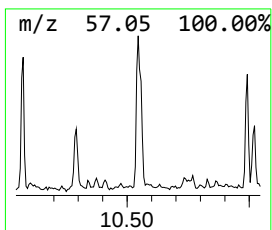
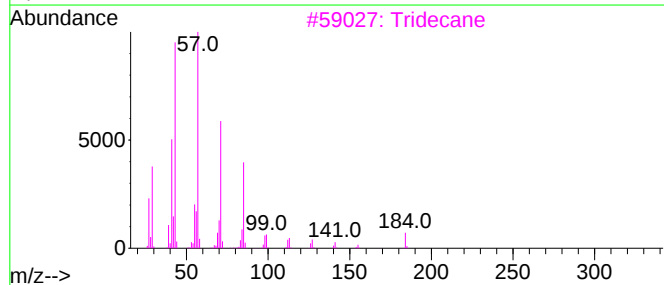
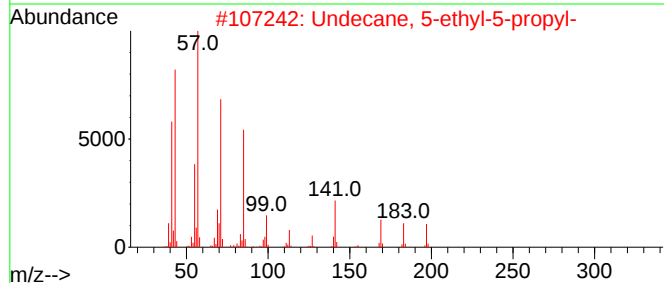
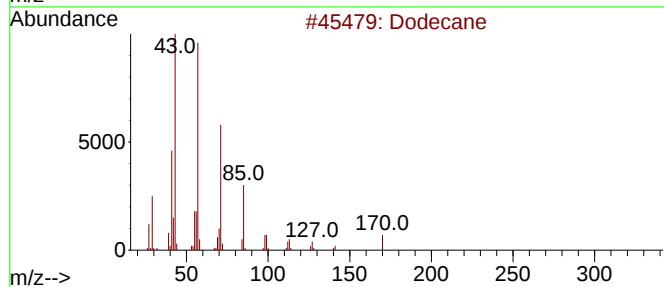
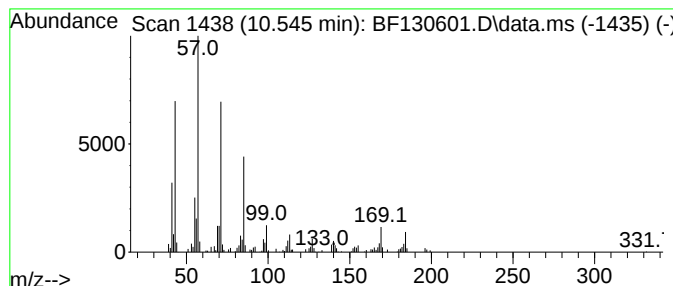
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Dodecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.545	4.00 ng	139371	Phenanthrene-d10	11.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	89
2			Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	83
3			Tridecane	184	C13H28	000629-50-5	81
4			Octadecane	254	C18H38	000593-45-3	74
5			Dodecane, 3-methyl-	184	C13H28	017312-57-1	74



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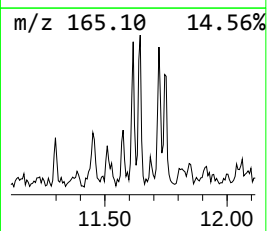
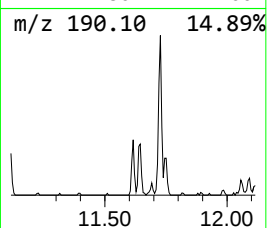
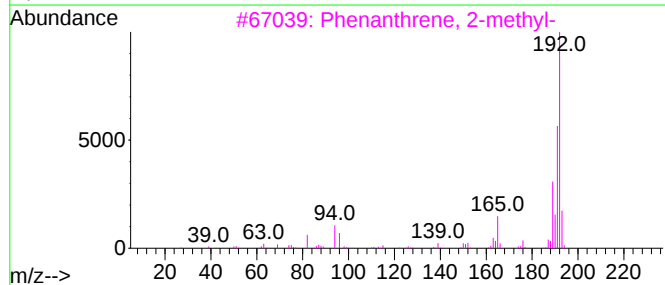
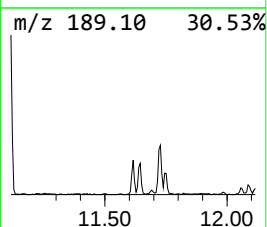
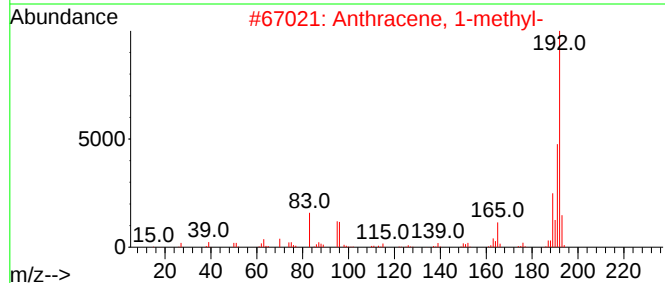
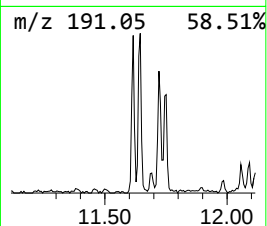
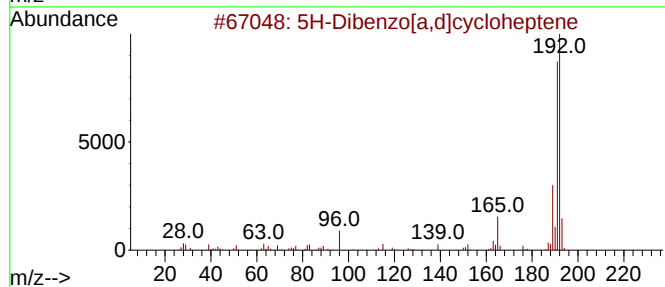
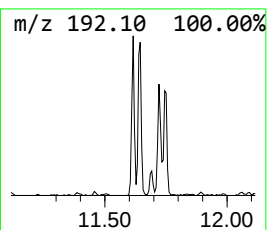
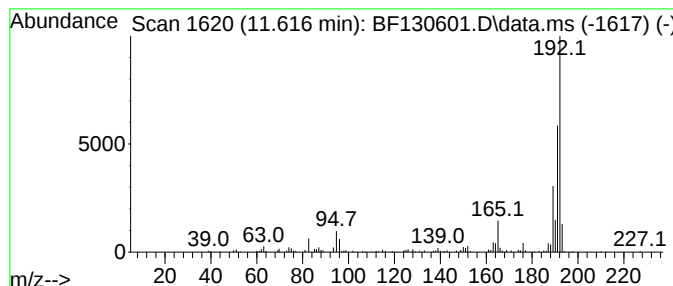
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Peak Number 3 5H-Dibenzo[a,d]cycloheptene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	2.55 ng	88795	Phenanthrene-d10	11.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	94
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91



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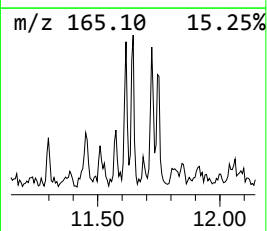
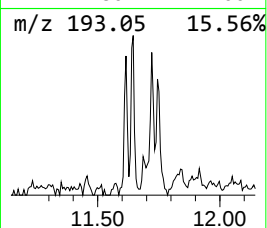
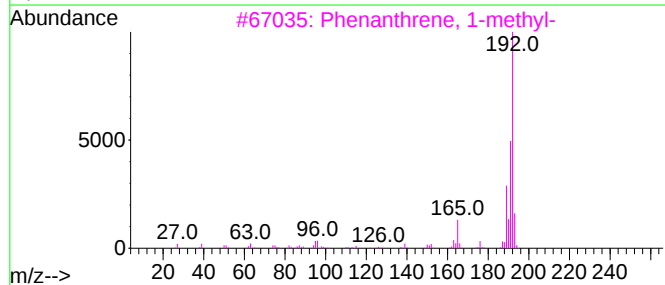
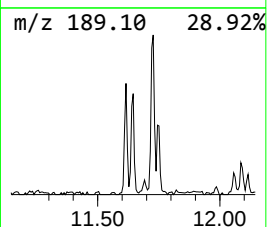
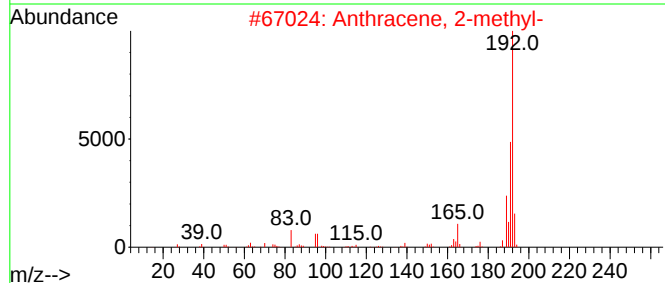
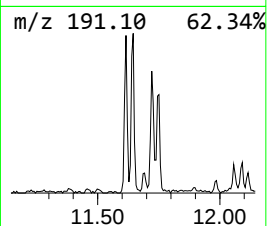
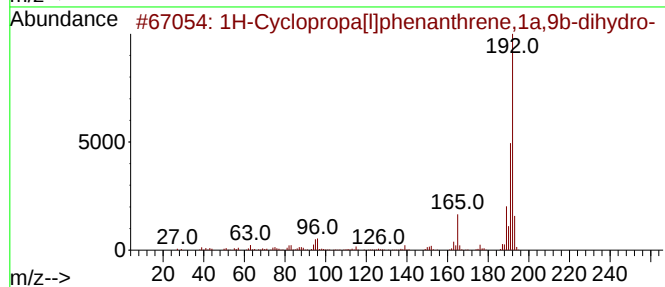
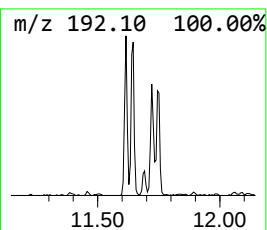
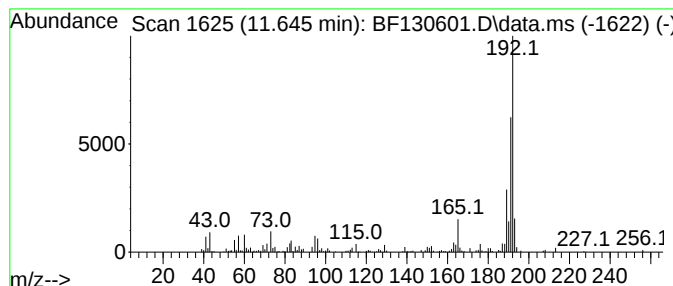
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1H-Cyclopropa[l]phenanthren... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.645	6.57 ng	229240	Phenanthrene-d10	11.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	98
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
Operator : CG\JU
Sample : N5046-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
R-2

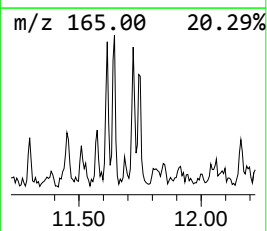
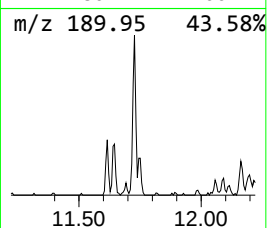
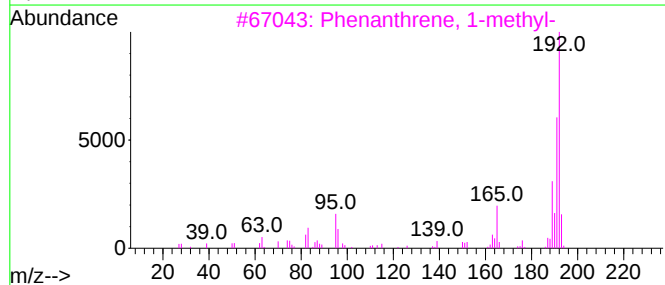
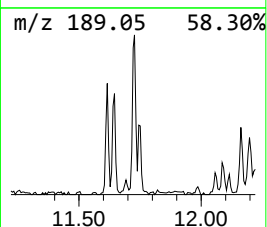
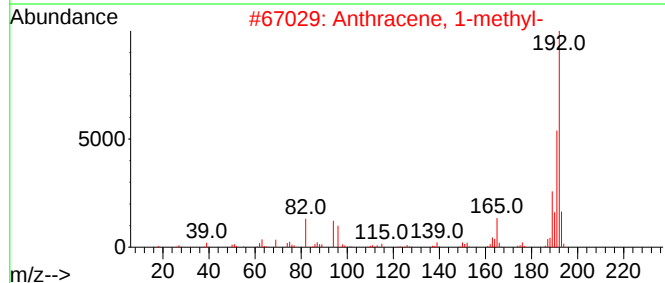
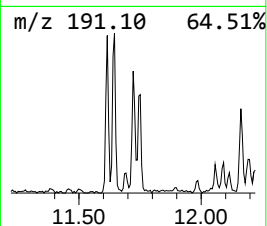
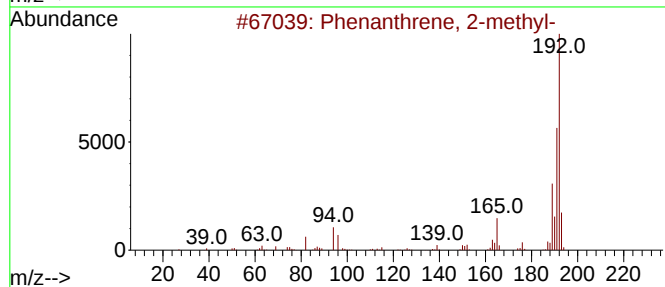
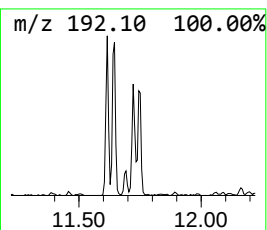
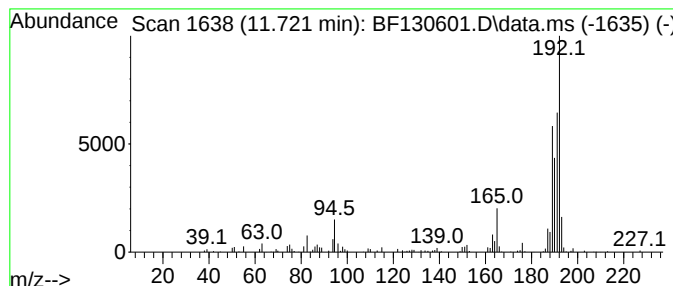
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.721	3.87 ng	134878	Phenanthrene-d10	11.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	78
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	70
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	60
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
Operator : CG\JU
Sample : N5046-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

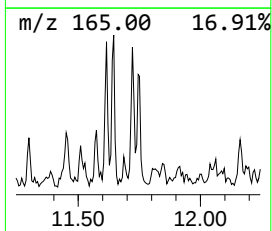
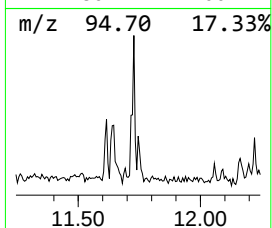
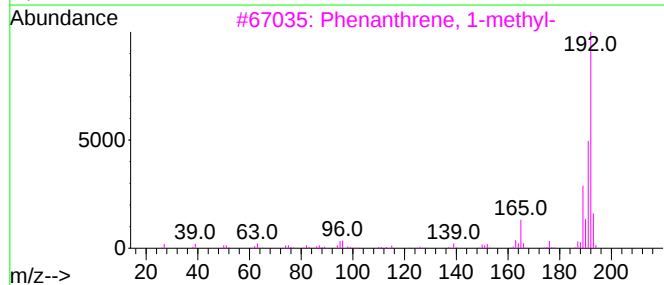
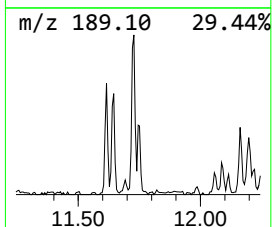
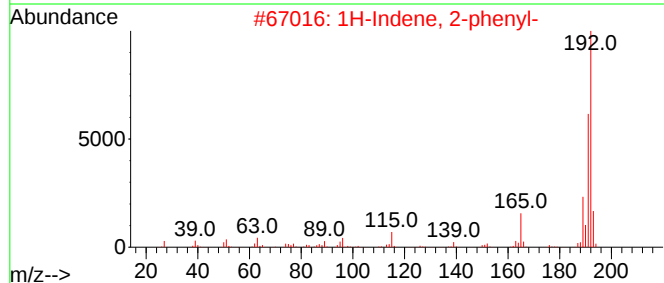
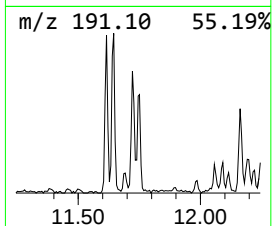
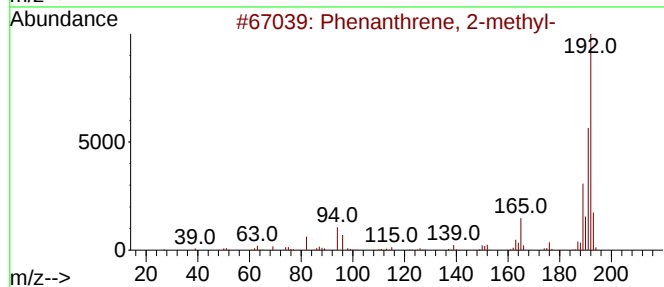
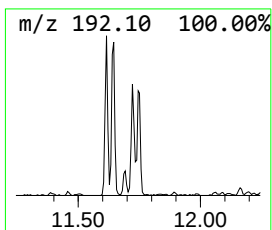
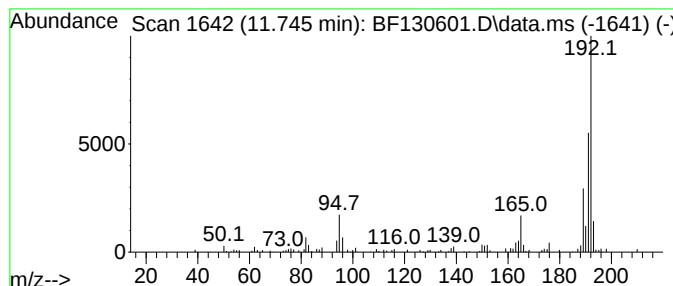
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BNA_F
ClientSampleId :
R-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.745	2.07 ng	72250	Phenanthrene-d10		11.116	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-		192	C15H12	002531-84-2	93
2	1H-Indene, 2-phenyl-		192	C15H12	004505-48-0	93
3	Phenanthrene, 1-methyl-		192	C15H12	000832-69-9	93
4	Phenanthrene, 4-methyl-		192	C15H12	000832-64-4	93
5	Anthracene, 2-methyl-		192	C15H12	000613-12-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
Operator : CG\JU
Sample : N5046-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
R-2

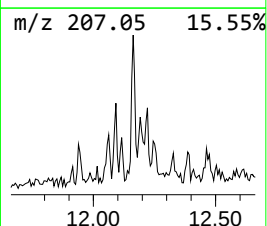
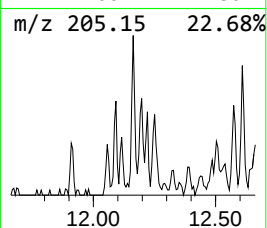
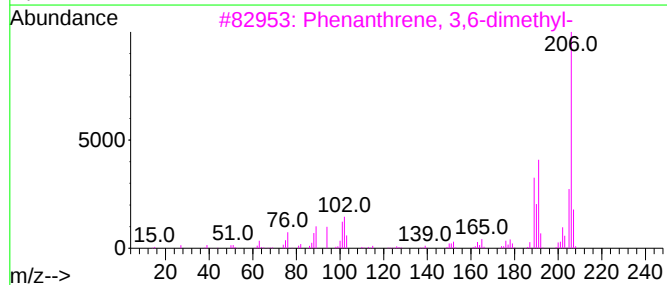
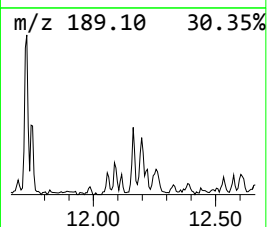
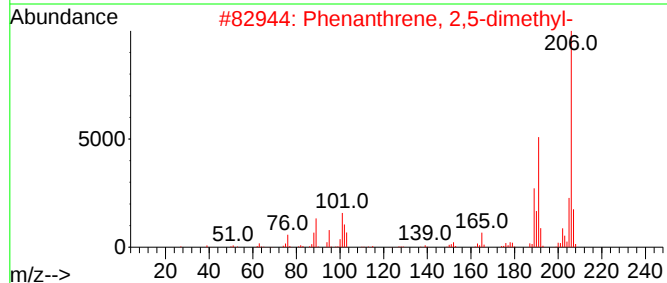
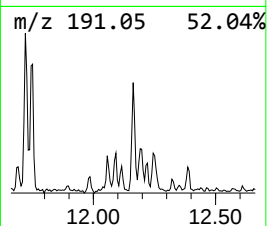
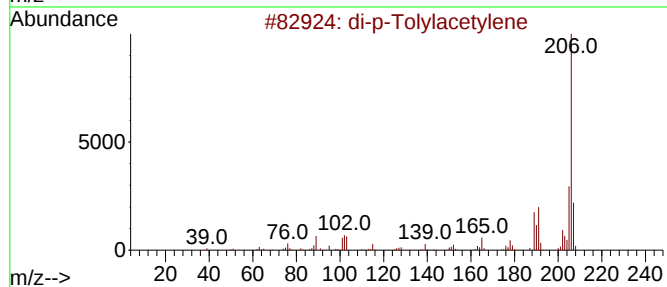
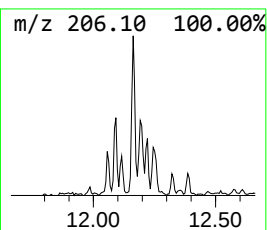
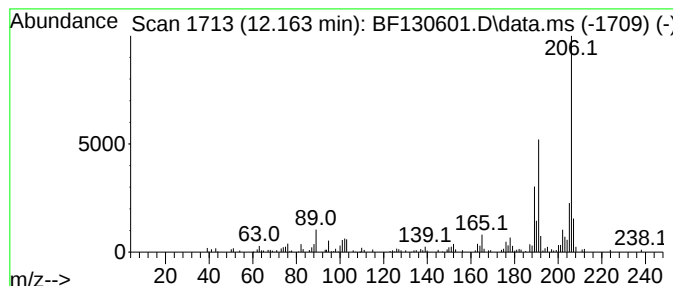
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 di-p-Tolylacetylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.163	2.59 ng	90416	Phenanthrene-d10	11.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			di-p-Tolylacetylene	206	C16H14	002789-88-0	93
2			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
3			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
4			1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	91
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
Operator : CG\JU
Sample : N5046-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
R-2

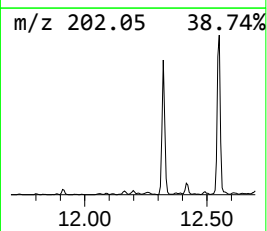
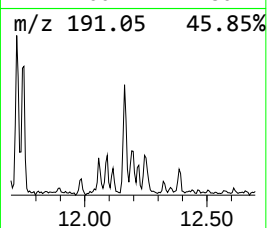
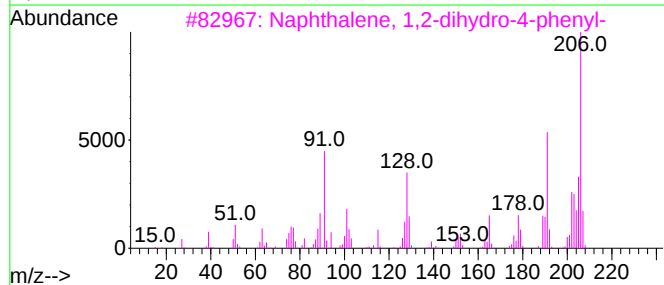
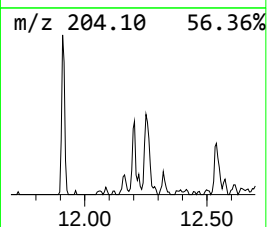
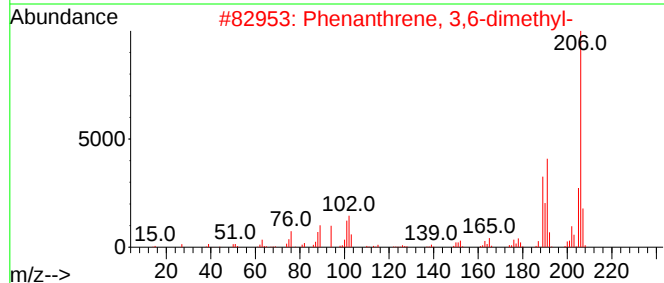
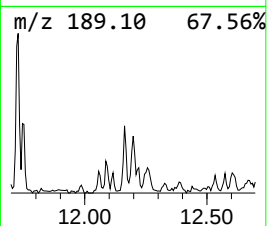
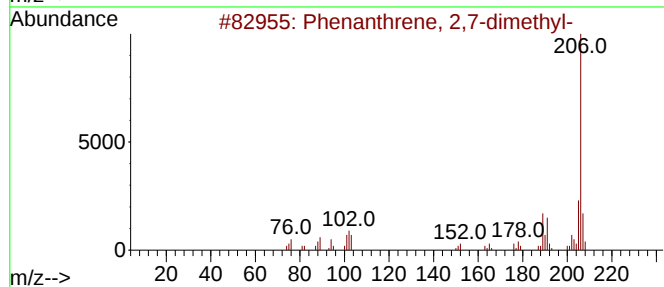
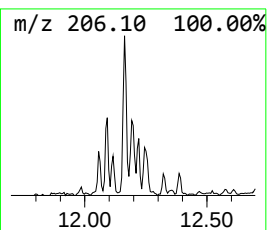
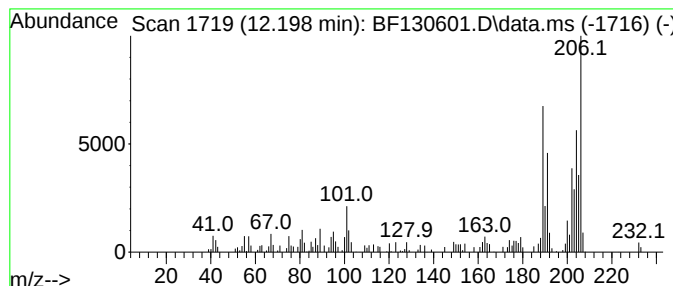
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2,7-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.198	2.33 ng	81125	Phenanthrene-d10	11.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	62
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	60
3			Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	46
4			Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	45
5			3,5-Dichloro-4-hydroxybenzoic acid	206	C7H4Cl2O3	003336-41-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
Operator : CG\JU
Sample : N5046-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

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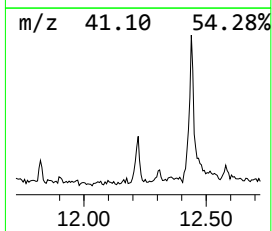
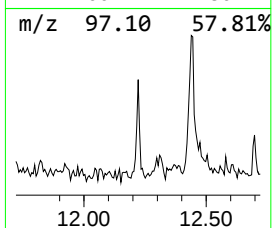
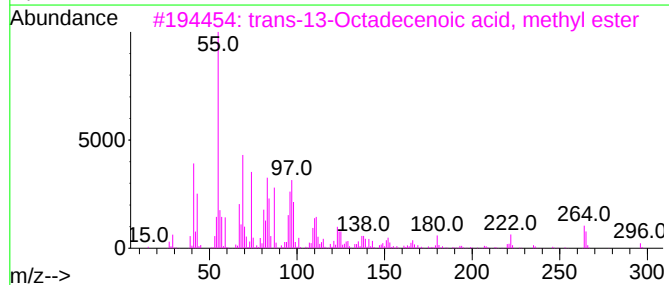
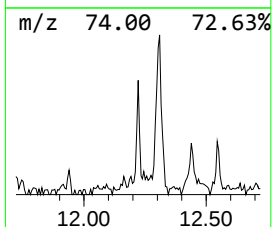
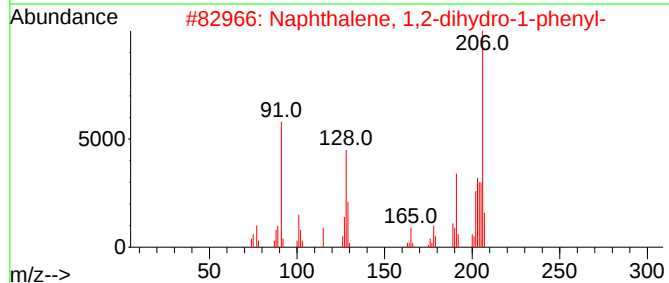
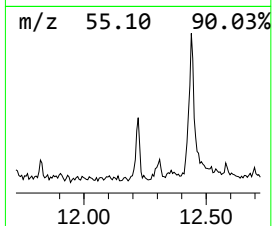
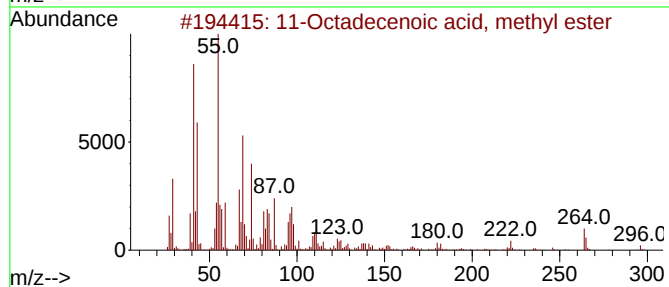
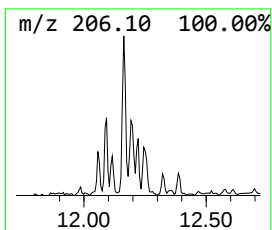
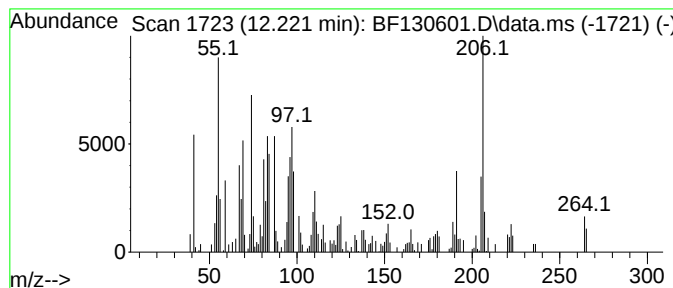
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11-Octadecenoic acid, methyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.221	2.62 ng	91458	Phenanthrene-d10	11.116

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	60
2		Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	55
3		trans-13-Octadecenoic acid, methyl ester	296	C19H36O2	1000333-61-3	55
4		1H-Fluorene, dodecahydro-	178	C13H22	005744-03-6	53
5		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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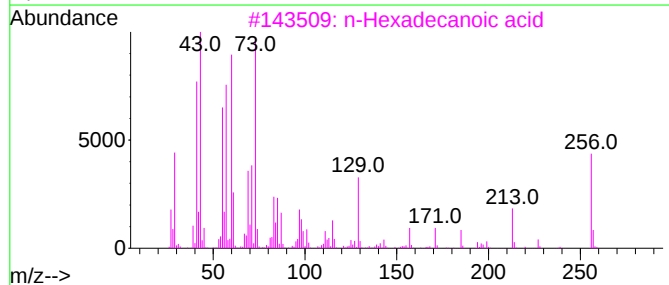
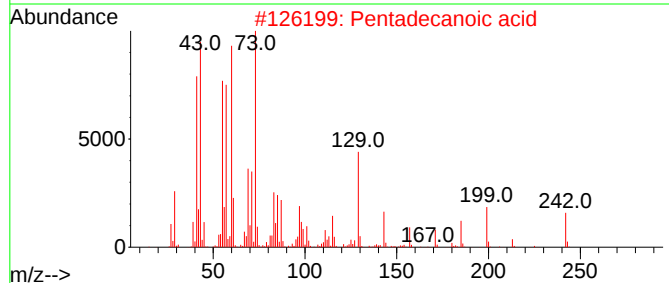
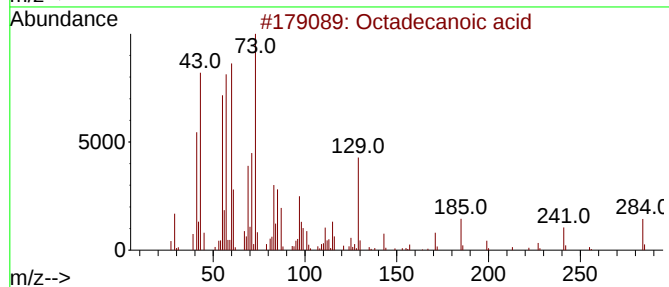
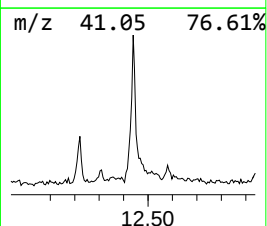
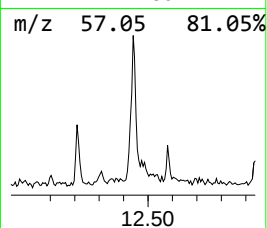
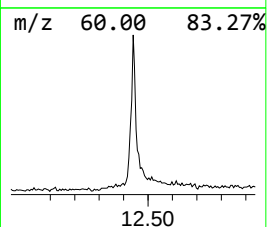
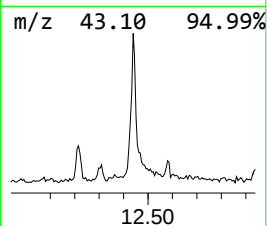
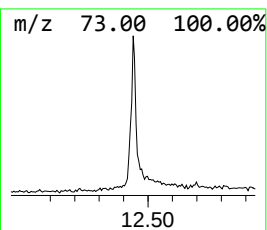
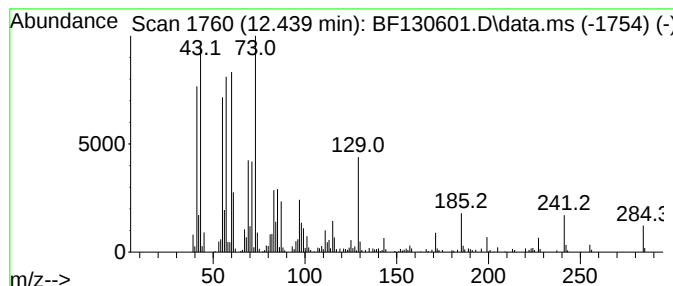
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.439	12.18 ng	405705	Chrysene-d12	13.745

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
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Operator : CG\JU
Sample : N5046-02
Misc :
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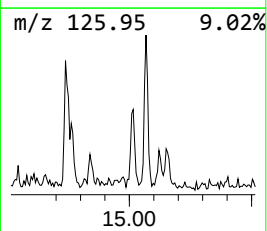
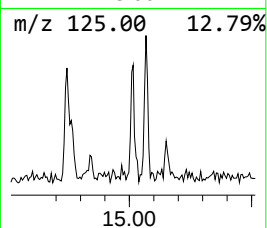
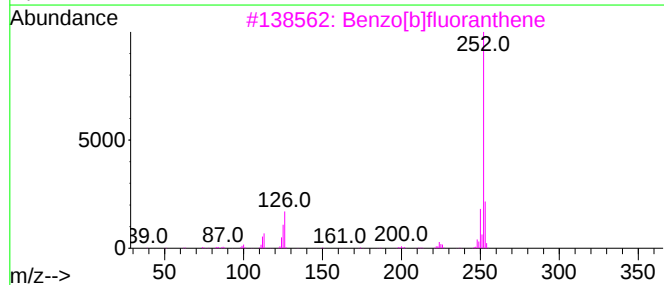
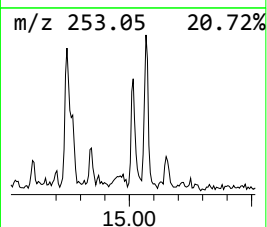
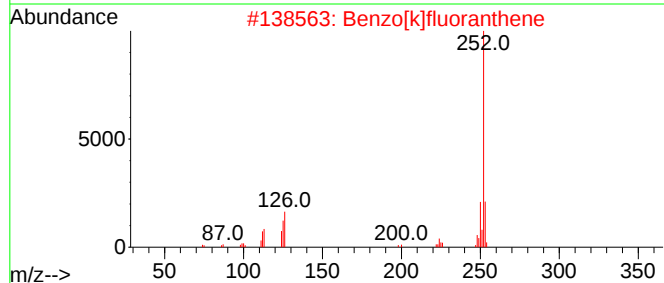
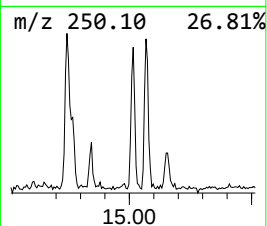
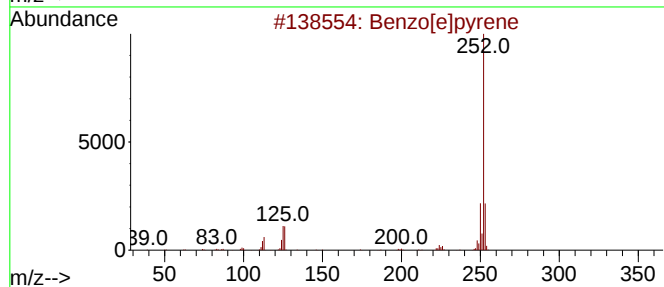
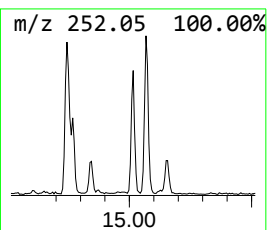
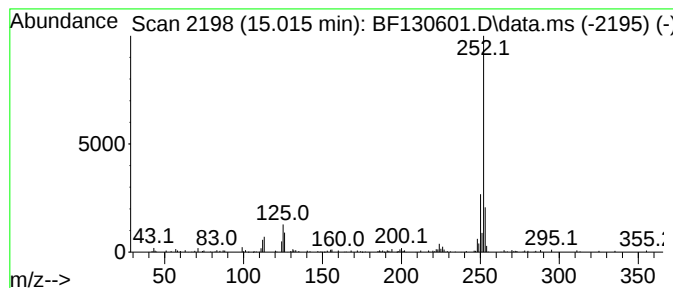
Instrument :
BNA_F
ClientSampleId :
R-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.015	2.27 ng	65838	Perylene-d12		15.127	
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[a]pyrene		252	C20H12	000050-32-8	93
5	Perylene		252	C20H12	000198-55-0	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
Operator : CG\JU
Sample : N5046-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
R-2

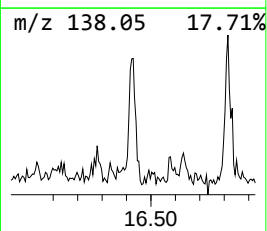
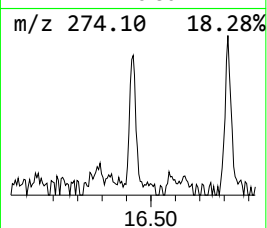
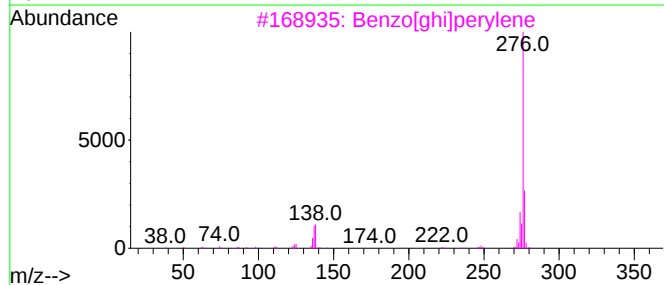
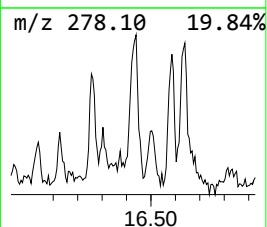
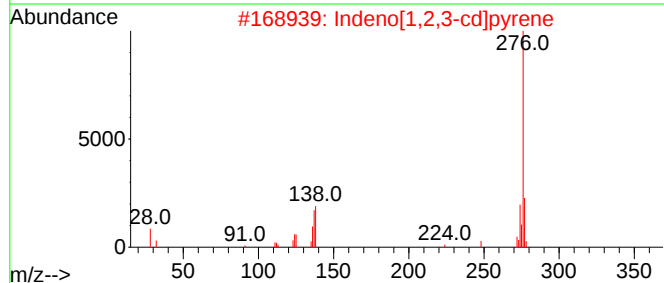
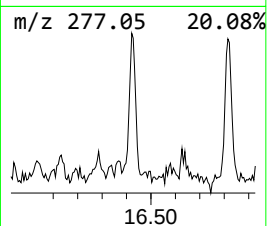
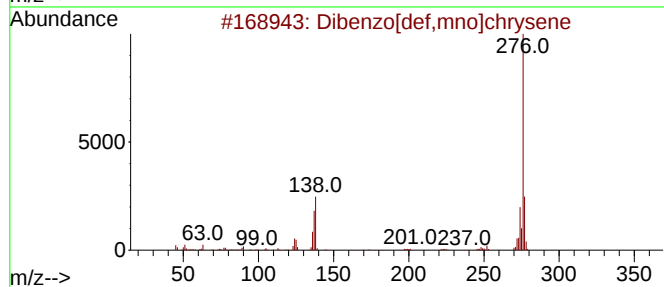
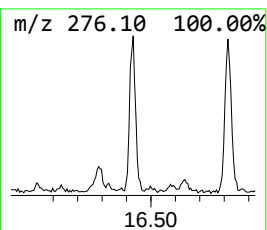
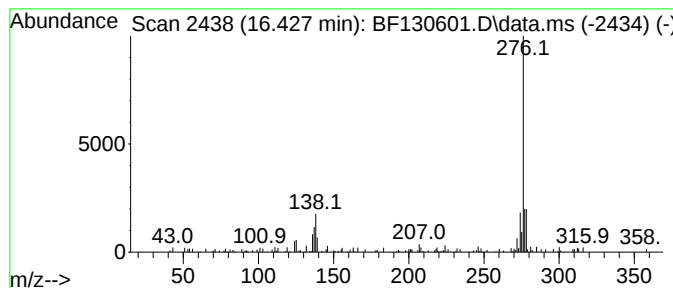
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,3,5-Triazin-2-amine, N-is... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.427	2.37 ng	68625	Perylene-d12	15.127

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	81
2		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	68
3		Benzo[ghi]perylene	276	C22H12	000191-24-2	68
4		1,3,5-Triazin-2-amine, N-isoprop...	276	C14H24N6	1000276-80-0	50
5		1,4-benzenediamine, N1,N1-dimeth...	276	C14H16N2O2S	1000401-83-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101022\
Data File : BF130601.D
Acq On : 10 Oct 2022 14:10
Operator : CG\JU
Sample : N5046-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
R-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.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
2-Pentanone, 4-...	4.781	20.3	ng	599456	1	6.604	589723	20.0
Dodecane	10.545	4.0	ng	139371	4	11.116	697546	20.0
5H-Dibenzo[a,d]...	11.616	2.5	ng	88795	4	11.116	697546	20.0
1H-Cyclopropa[l...	11.645	6.6	ng	229240	4	11.116	697546	20.0
Anthracene, 1-m...	11.721	3.9	ng	134878	4	11.116	697546	20.0
Phenanthrene, 2...	11.745	2.1	ng	72250	4	11.116	697546	20.0
di-p-Tolylacety...	12.163	2.6	ng	90416	4	11.116	697546	20.0
Phenanthrene, 2...	12.198	2.3	ng	81125	4	11.116	697546	20.0
11-Octadecenoic...	12.221	2.6	ng	91458	4	11.116	697546	20.0
Octadecanoic acid	12.439	12.2	ng	405705	5	13.745	666215	20.0
Benzo[e]pyrene	15.015	2.3	ng	65838	6	15.127	579232	20.0
1,3,5-Triazin-2...	16.427	2.4	ng	68625	6	15.127	579232	20.0