

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
 Data File : BF131977.D
 Acq On : 09 Jan 2023 22:39
 Operator : CG\JU
 Sample : 01065-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-210

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-BF010923.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131977.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.128	3	7	16	rVB	44968	57397	2.51%	0.280%
2	4.987	488	493	506	rVB	580424	805168	35.15%	3.923%
3	5.398	559	563	573	rBV	1922216	1884680	82.29%	9.182%
4	6.422	732	737	740	rBV	1866964	1914738	83.60%	9.328%
5	6.781	795	798	802	rBV	523999	428791	18.72%	2.089%
6	7.357	890	896	899	rBV	1290699	1242601	54.25%	6.054%
7	8.075	1013	1018	1021	rBV	601837	550391	24.03%	2.681%
8	9.157	1196	1202	1205	rBV	2588318	2290370	100.00%	11.158%
9	9.833	1313	1317	1320	rBV	802230	639378	27.92%	3.115%
10	9.863	1320	1322	1326	rVB	61911	47250	2.06%	0.230%
11	10.380	1407	1410	1416	rVB	43950	38033	1.66%	0.185%
12	10.551	1435	1439	1443	rVB	160263	130871	5.71%	0.638%
13	10.627	1448	1452	1455	rBV	1819265	1485179	64.84%	7.235%
14	11.222	1548	1553	1557	rVB	27353	26076	1.14%	0.127%
15	11.327	1567	1571	1573	rBV	738563	708944	30.95%	3.454%
16	11.351	1573	1575	1578	rVV	517526	405849	17.72%	1.977%
17	11.398	1578	1583	1586	rVV	103938	97600	4.26%	0.475%
18	11.557	1604	1610	1617	rBV2	49103	60600	2.65%	0.295%
19	11.827	1651	1656	1658	rBV	72312	62026	2.71%	0.302%
20	11.857	1658	1661	1664	rVV	107972	95122	4.15%	0.463%
21	11.904	1664	1669	1672	rVV2	26354	31383	1.37%	0.153%
22	11.939	1672	1675	1677	rVV	117040	111280	4.86%	0.542%
23	11.963	1677	1679	1683	rVB	45658	37458	1.64%	0.182%
24	12.121	1704	1706	1709	rVV	44862	38209	1.67%	0.186%
25	12.157	1709	1712	1715	rVB	40558	37479	1.64%	0.183%
26	12.380	1746	1750	1752	rBV	39629	45468	1.99%	0.222%
27	12.469	1762	1765	1769	rVB2	37924	40076	1.75%	0.195%
28	12.539	1774	1777	1781	rBV	737137	691054	30.17%	3.367%
29	12.692	1796	1803	1806	rBV4	25212	38335	1.67%	0.187%
30	12.774	1810	1817	1822	rVV2	564541	668673	29.19%	3.258%
31	12.821	1822	1825	1829	rVB2	37307	41063	1.79%	0.200%
32	12.916	1837	1841	1845	rVB	2153703	2064685	90.15%	10.058%
33	12.998	1848	1855	1858	rBV2	41520	76624	3.35%	0.373%
34	13.080	1865	1869	1870	rBV2	37596	42885	1.87%	0.209%
35	13.104	1870	1873	1875	rVB	94333	87748	3.83%	0.427%
36	13.168	1881	1884	1887	rBV	75388	57829	2.52%	0.282%

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Integrator: RTE

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Filtering: 5

Sampling : 1

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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-BF010923.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.204	1887	1890	1892	rVV2	63728	65232	2.85%	0.318%
38	13.227	1892	1894	1898	rVB	31317	28336	1.24%	0.138%
39	13.298	1903	1906	1909	rBV	34641	28607	1.25%	0.139%
40	13.327	1909	1911	1915	rBV4	29429	30993	1.35%	0.151%
41	13.615	1957	1960	1964	rVB	40198	40890	1.79%	0.199%
42	13.721	1973	1978	1981	rBV	59873	64304	2.81%	0.313%
43	13.751	1981	1983	1985	rVV	41139	31137	1.36%	0.152%
44	13.815	1991	1994	2003	rVB3	163824	245994	10.74%	1.198%
45	13.974	2013	2021	2024	rBV2	550616	760406	33.20%	3.704%
46	13.998	2024	2025	2032	rVB	238479	206747	9.03%	1.007%
47	14.080	2037	2039	2041	rVB	42366	29251	1.28%	0.143%
48	14.133	2044	2048	2053	rBV8	19609	38161	1.67%	0.186%
49	14.363	2082	2087	2090	rBV	51077	48239	2.11%	0.235%
50	14.398	2090	2093	2097	rVB3	23159	27362	1.19%	0.133%
51	14.457	2097	2103	2106	rBV5	32090	59735	2.61%	0.291%
52	14.680	2136	2141	2143	rBV3	20609	26881	1.17%	0.131%
53	15.015	2193	2198	2201	rBV	220580	330615	14.44%	1.611%
54	15.121	2212	2216	2220	rVB	49663	63346	2.77%	0.309%
55	15.315	2245	2249	2255	rVB	118032	153959	6.72%	0.750%
56	15.374	2255	2259	2265	rBV	170212	205659	8.98%	1.002%
57	15.439	2265	2270	2274	rBV	332441	421074	18.38%	2.051%
58	15.657	2303	2307	2315	rVB5	40100	61260	2.67%	0.298%
59	15.886	2340	2346	2349	rBV6	18050	32877	1.44%	0.160%
60	15.945	2352	2356	2361	rVV6	26642	58545	2.56%	0.285%
61	15.986	2361	2363	2370	rVB3	15479	27237	1.19%	0.133%
62	16.892	2511	2517	2524	rBV3	73952	145856	6.37%	0.711%
63	17.068	2543	2547	2552	rBV3	19446	38086	1.66%	0.186%
64	17.127	2554	2557	2563	rVB3	16908	29893	1.31%	0.146%
65	17.333	2586	2592	2602	rVB	53198	114809	5.01%	0.559%
66	17.580	2627	2634	2637	rBV	14582	30048	1.31%	0.146%

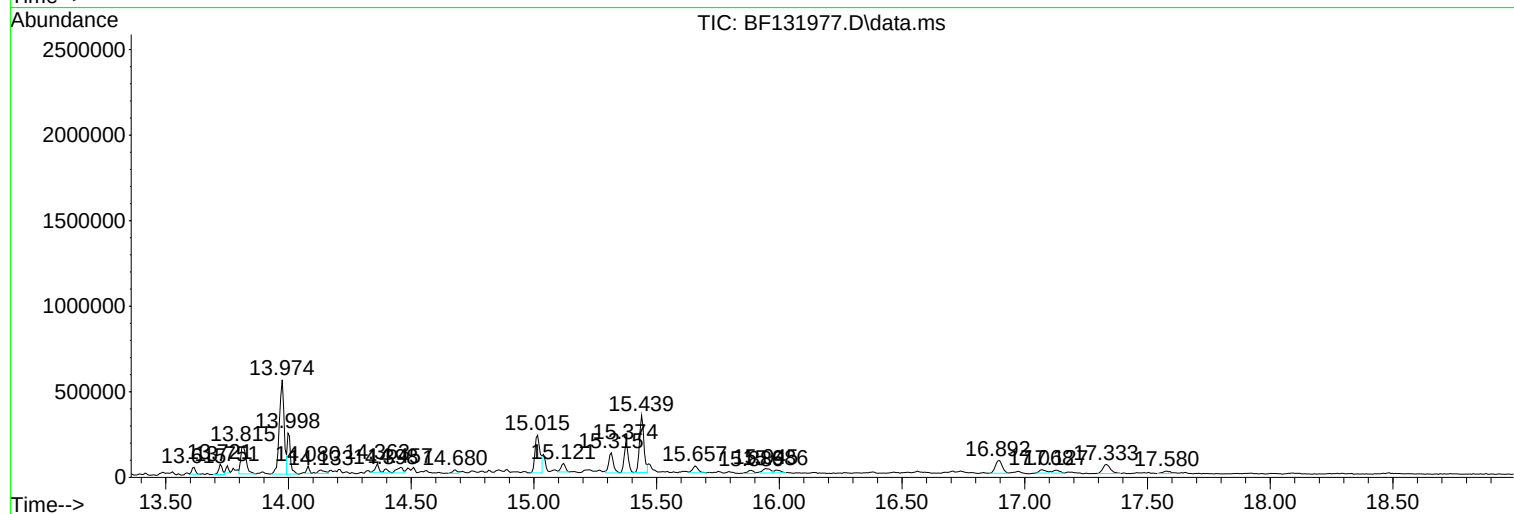
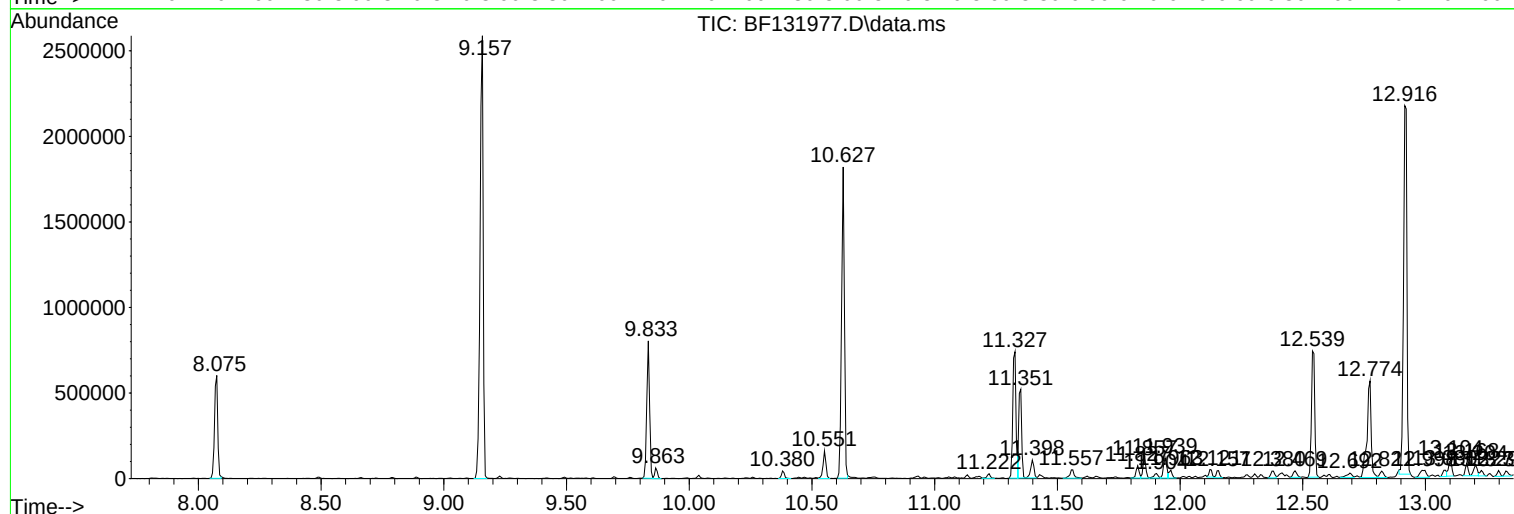
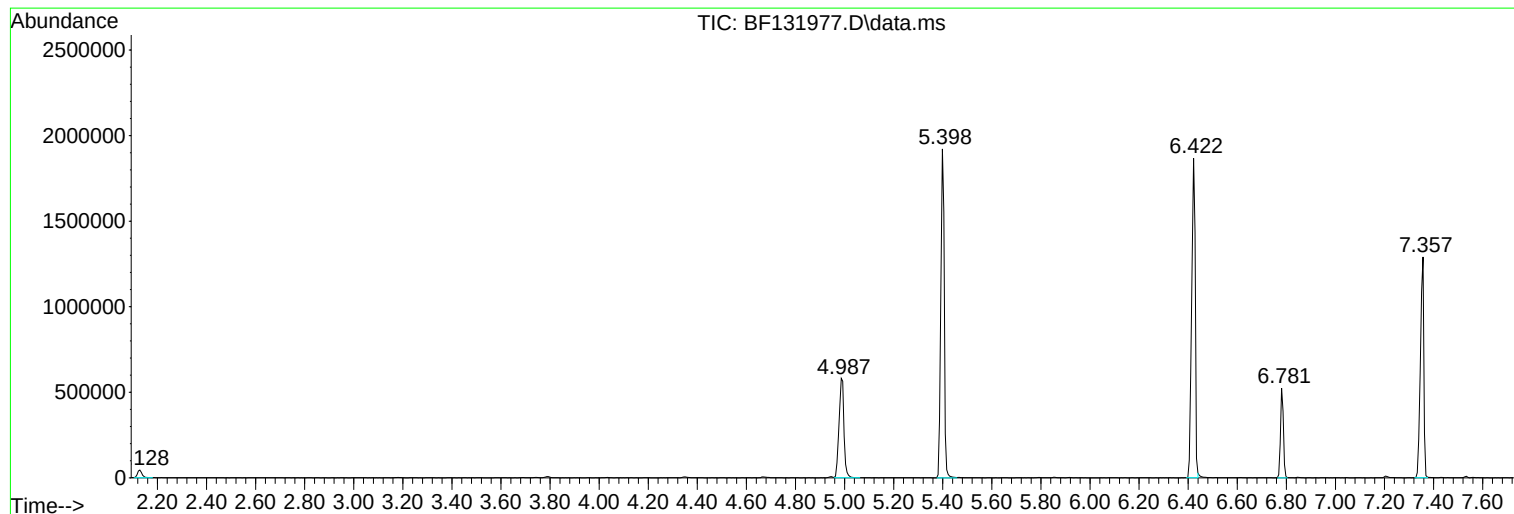
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Instrument :
BNA_F
ClientSampleId :
VNJ-210

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.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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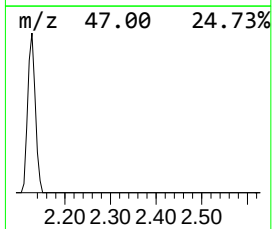
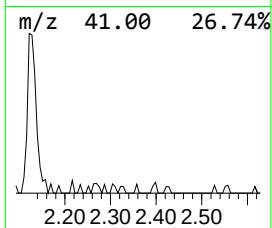
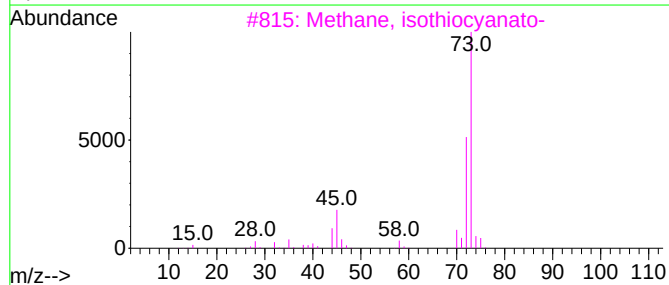
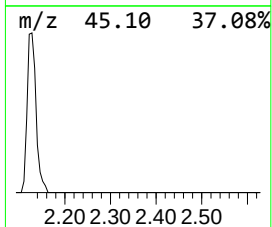
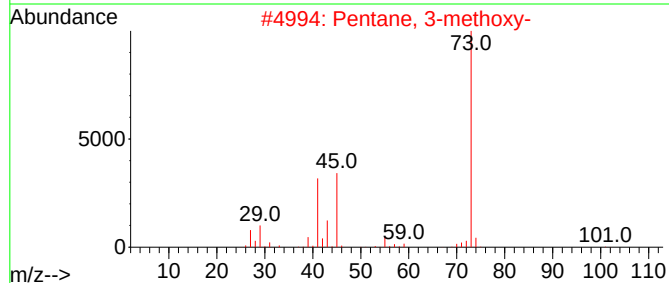
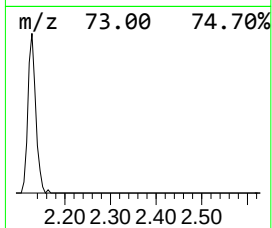
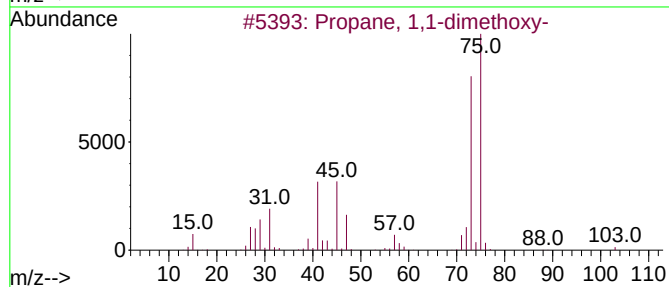
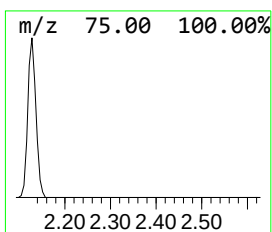
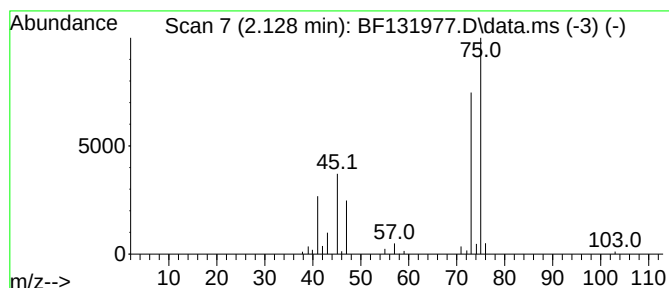
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.128	2.68 ng	57397	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	14
3			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9
4			Thiocyanic acid, methyl ester	73	C2H3NS	000556-64-9	9
5			1,3-Dioxolane, 2-(1-bromoethyl)-	180	C5H9BrO2	005267-73-2	9



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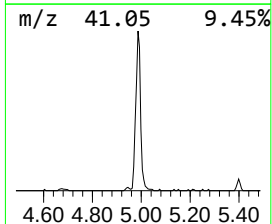
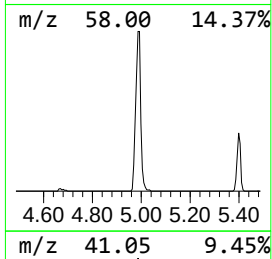
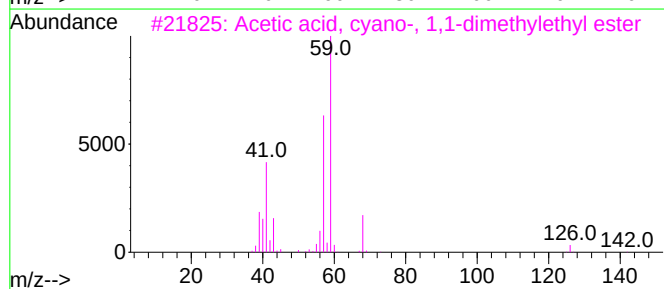
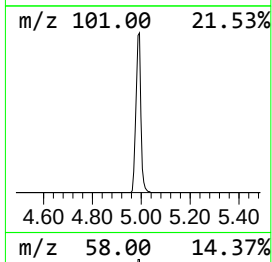
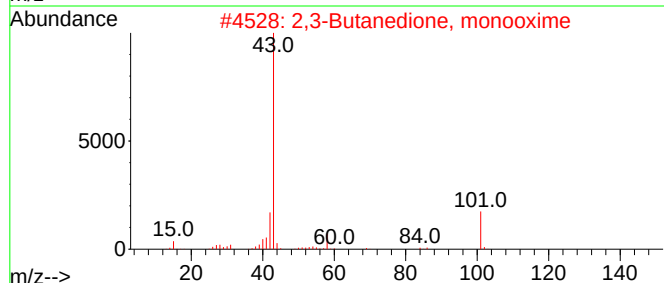
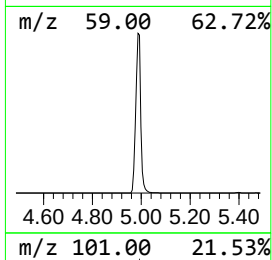
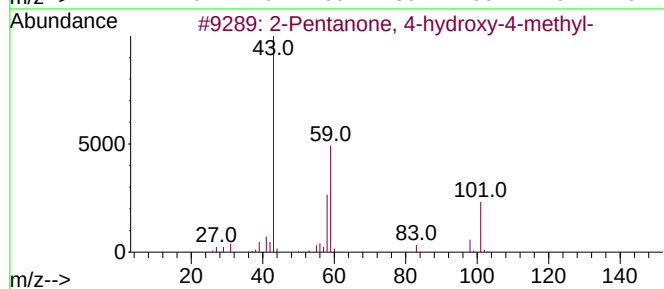
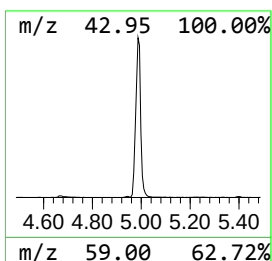
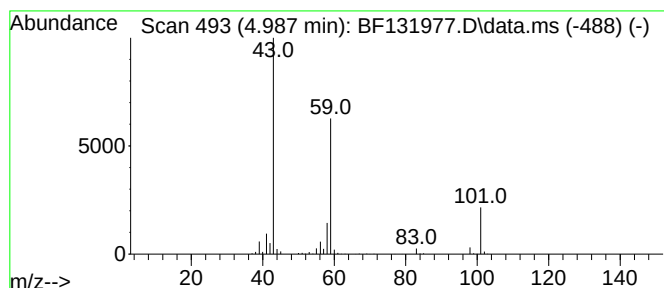
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.987	37.56 ng	805168	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



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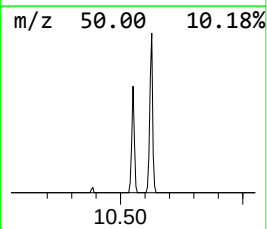
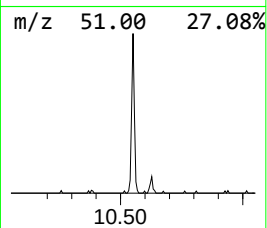
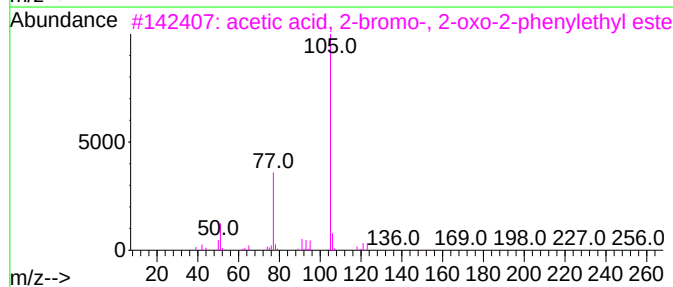
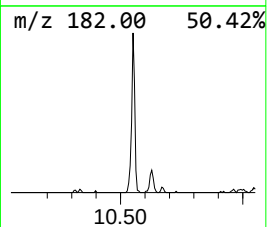
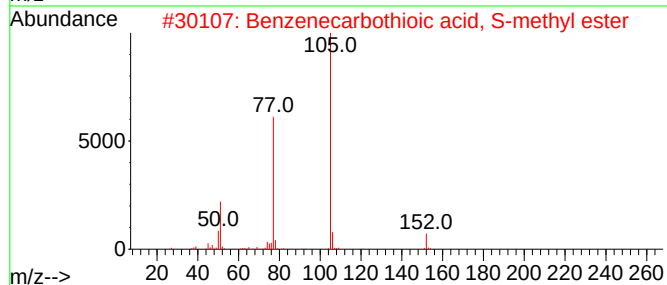
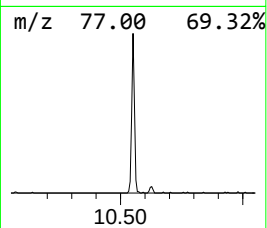
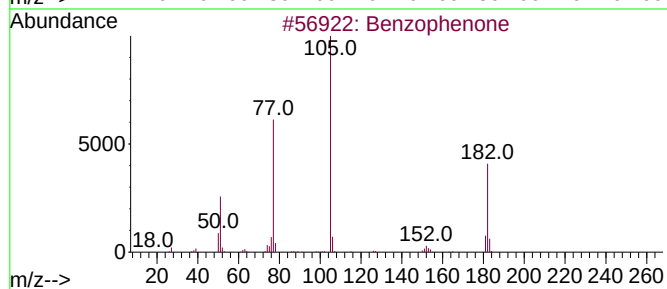
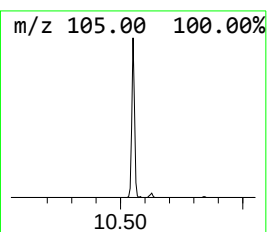
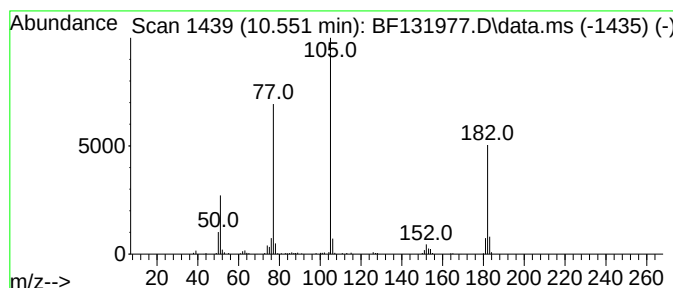
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.551	4.09 ng	130871	Acenaphthene-d10	9.833

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	52
3			acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	50
4			2-Phenyl-4-ethylidene-2-oxazolin...	187	C11H9NO2	037791-41-6	50
5			Benzoylformic acid	150	C8H6O3	000611-73-4	49



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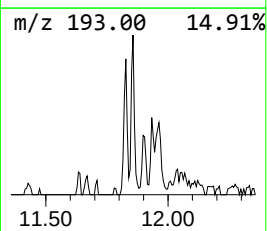
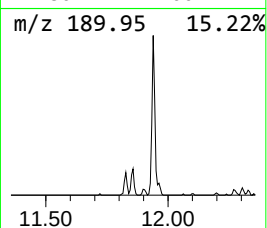
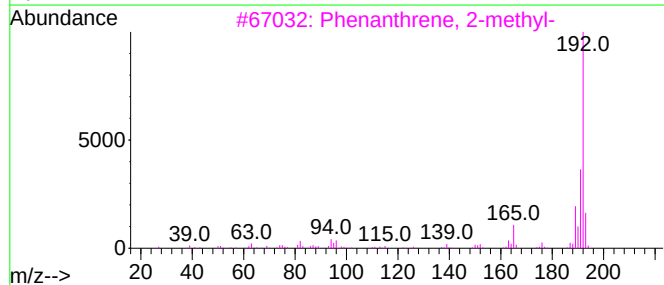
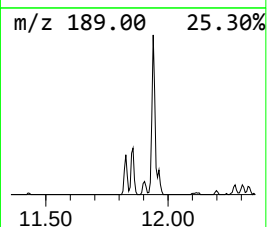
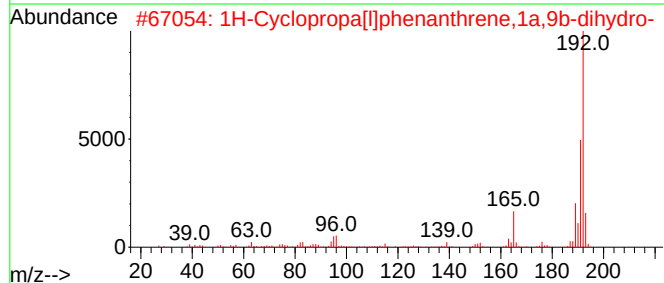
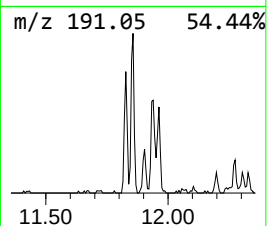
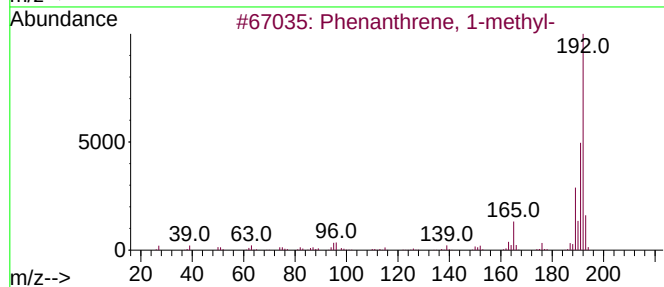
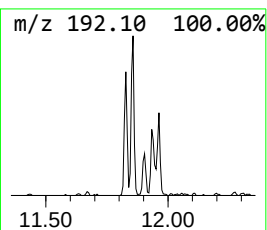
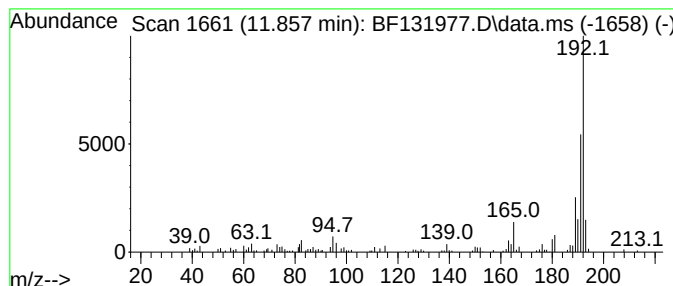
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.857	2.68 ng	95122	Phenanthrene-d10	11.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
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Acq On : 09 Jan 2023 22:39
Operator : CG\JU
Sample : 01065-01
Misc :
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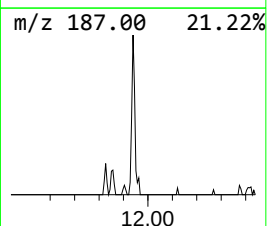
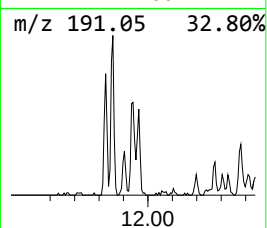
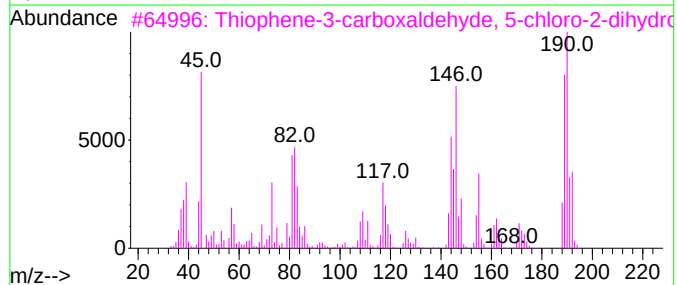
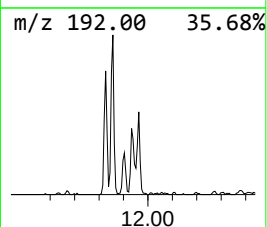
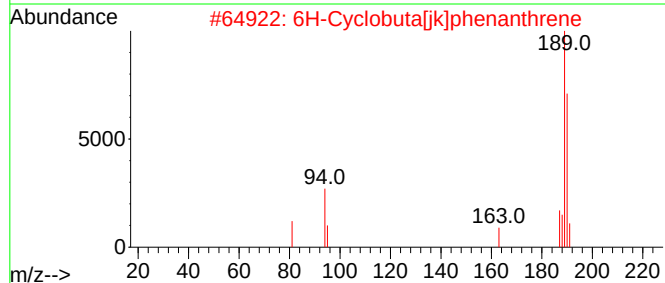
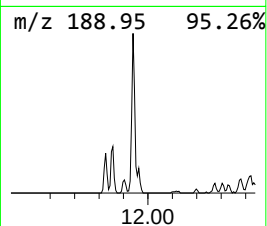
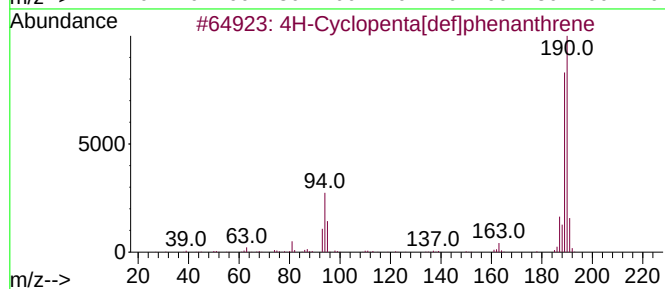
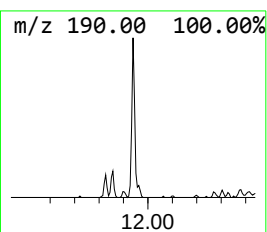
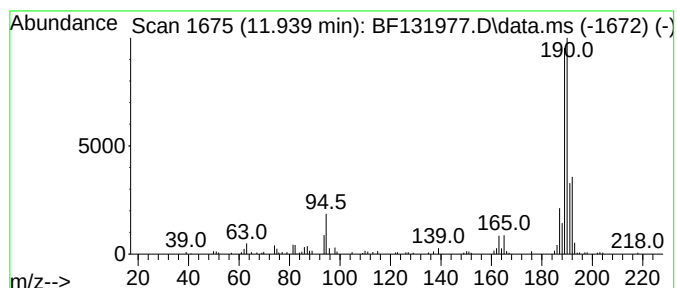
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.939	3.14 ng	111280	Phenanthrene-d10	11.327	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4	Methyl diselenide	190	C2H6Se2	007101-31-7	49
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
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Sample : 01065-01
Misc :
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Instrument :
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ClientSampleId :
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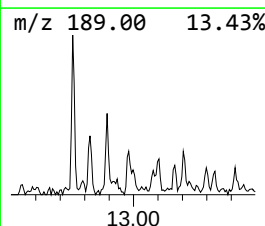
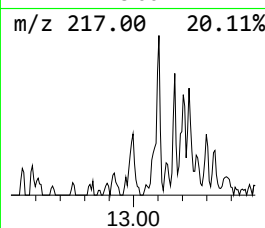
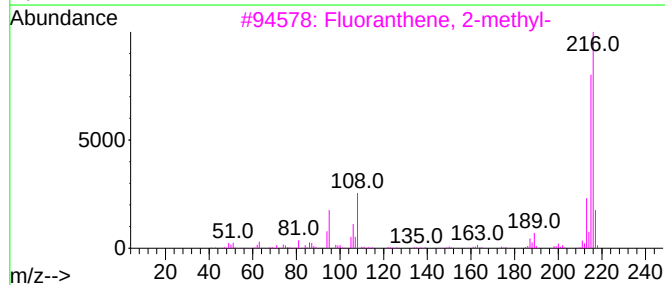
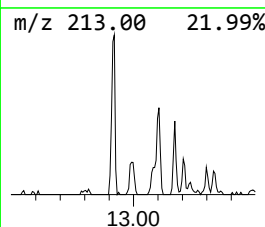
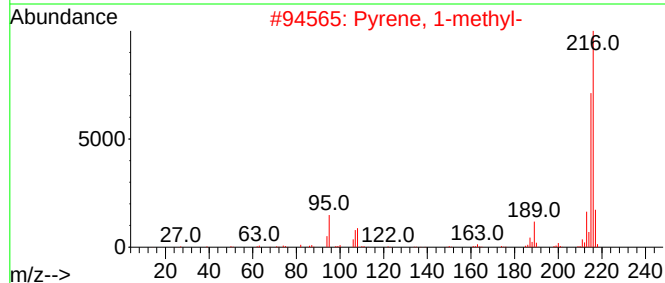
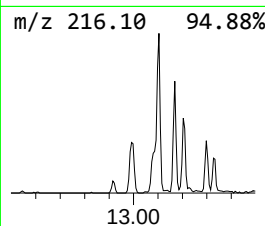
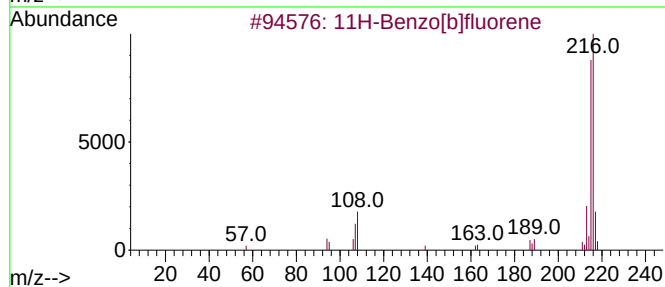
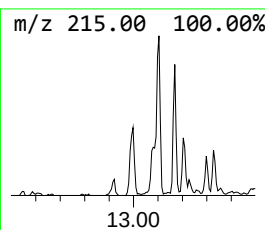
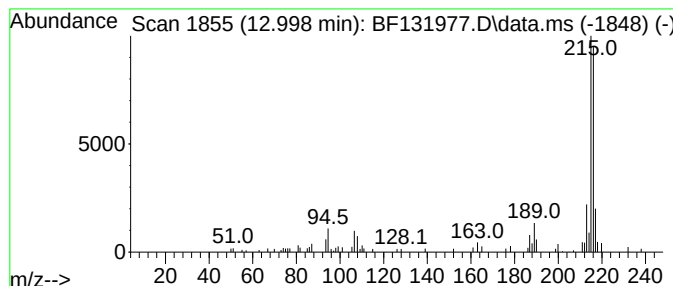
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.998	2.02 ng	76624	Chrysene-d12	13.974

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	74
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
Operator : CG\JU
Sample : 01065-01
Misc :
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Instrument :
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ClientSampleId :
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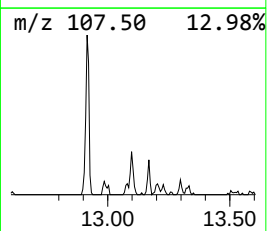
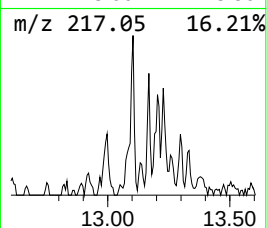
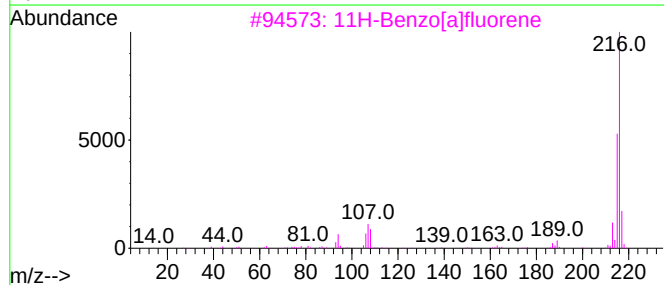
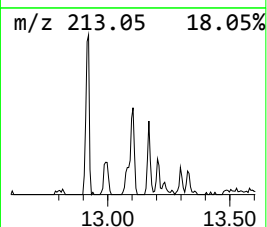
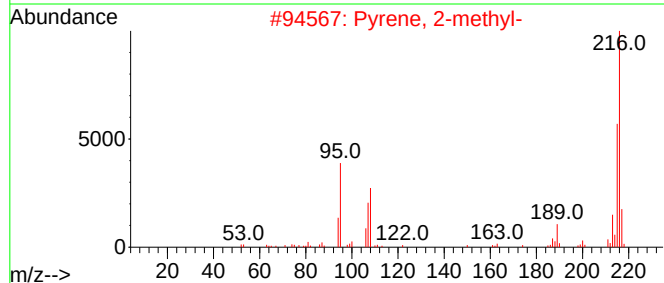
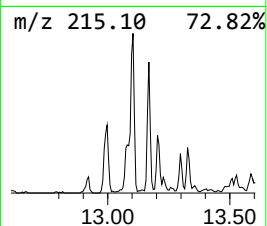
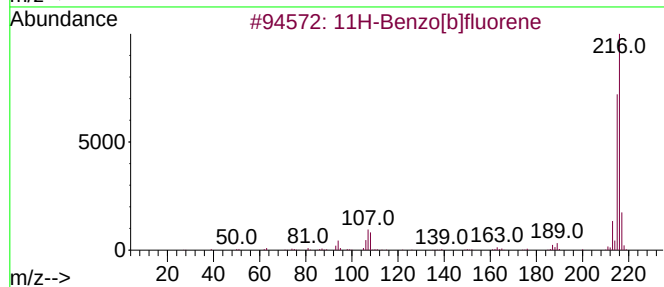
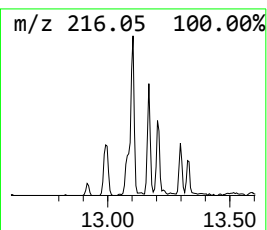
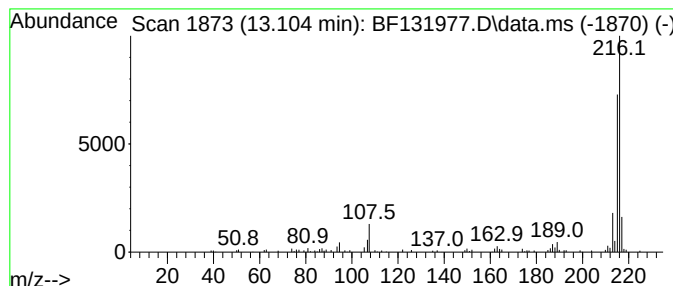
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pyrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.104	2.31 ng	87748	Chrysene-d12	13.974

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
4		Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
5		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
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Sample : 01065-01
Misc :
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Instrument :
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ClientSampleId :
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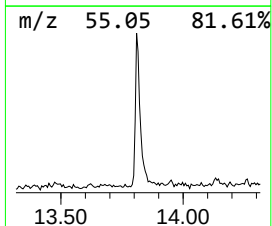
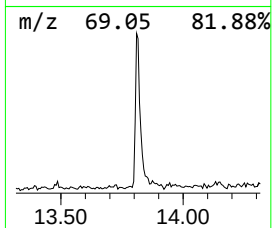
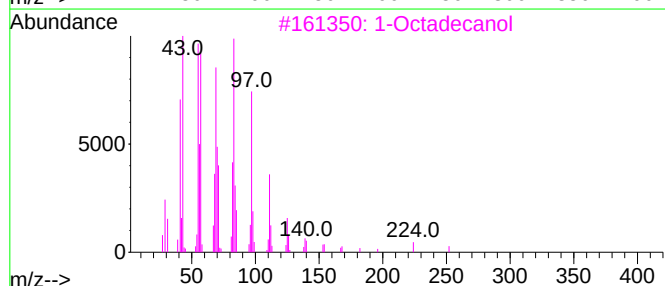
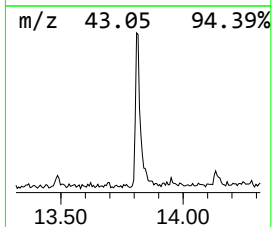
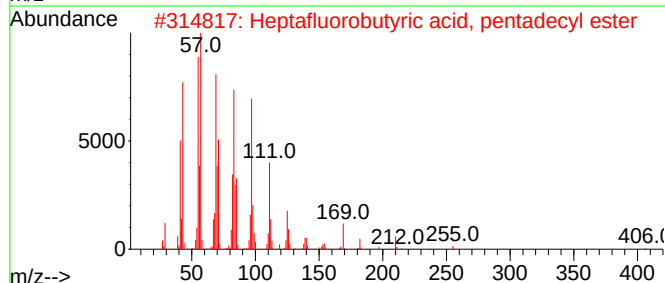
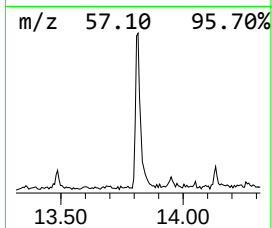
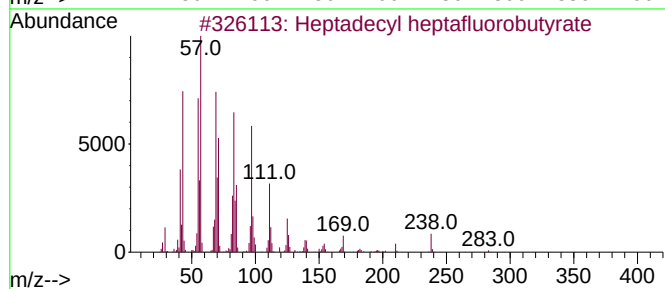
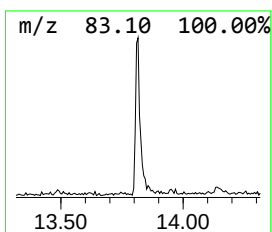
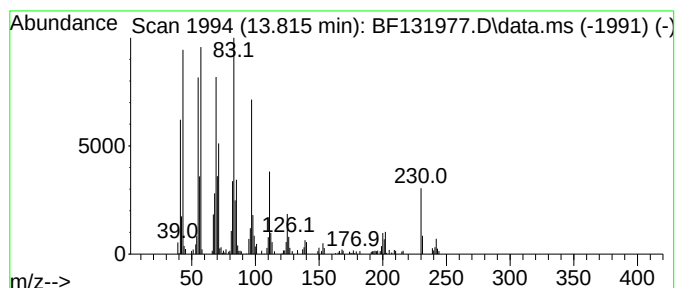
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecyl heptafluorobutyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.816	6.47 ng	245994	Chrysene-d12	13.974

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	89
2			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	89
3			1-Octadecanol	270	C18H38O	000112-92-5	81
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	76
5			1-Tetracosene	336	C24H48	010192-32-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
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Sample : 01065-01
Misc :
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Instrument :
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ClientSampleId :
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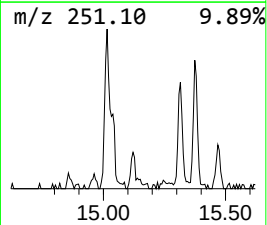
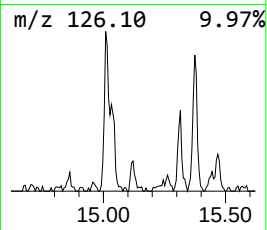
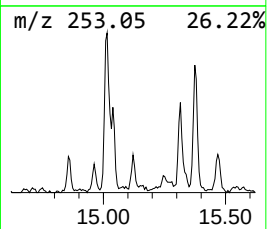
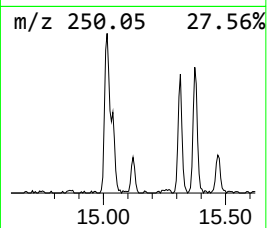
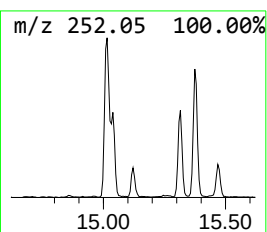
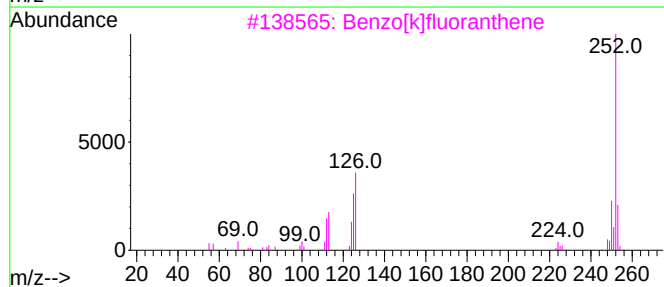
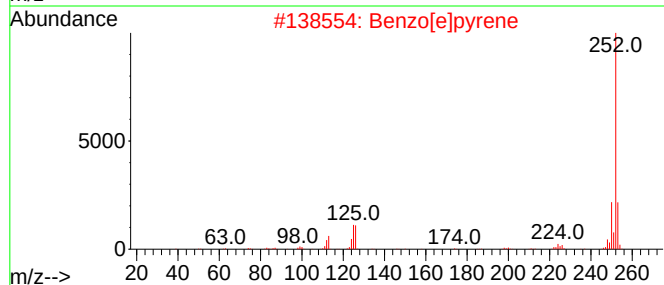
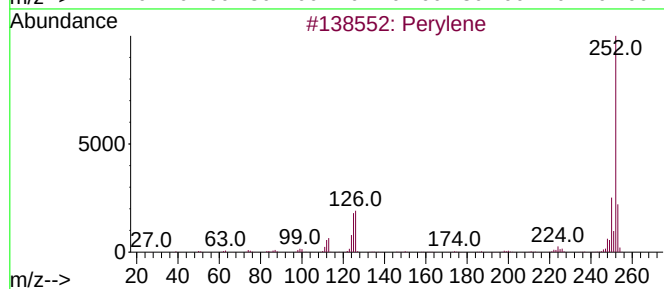
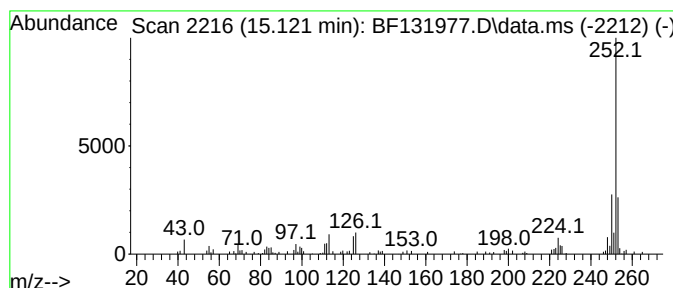
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Perylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.121	3.01 ng	63346	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	94
2			Benzo[e]pyrene	252	C20H12	000192-97-2	91
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	87
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	86
5			Benzo[a]pyrene	252	C20H12	000050-32-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
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Misc :
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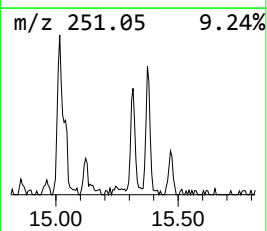
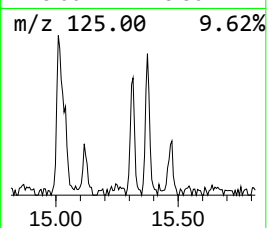
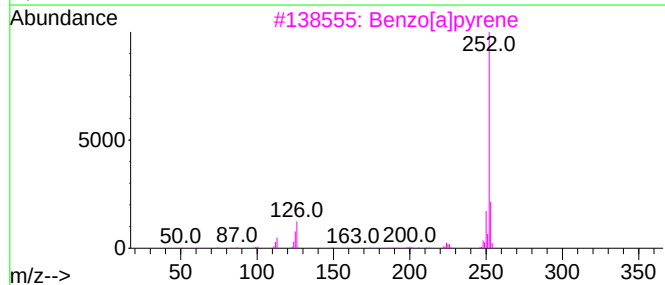
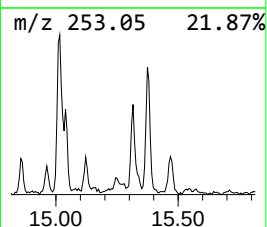
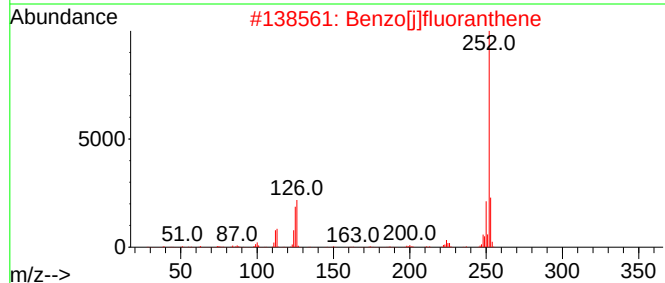
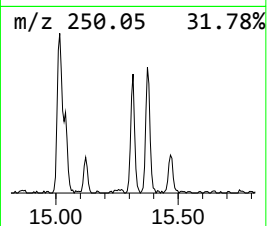
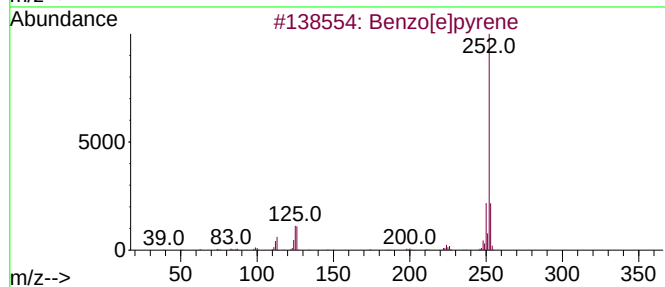
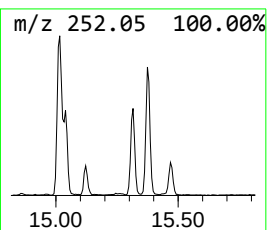
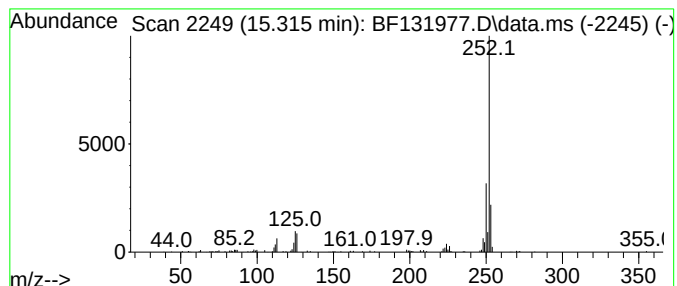
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.315	7.31 ng	153959	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	95
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			Benzo[b]fluoranthene	252	C20H12	000205-99-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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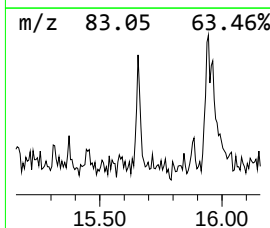
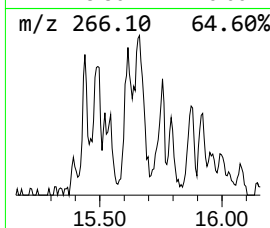
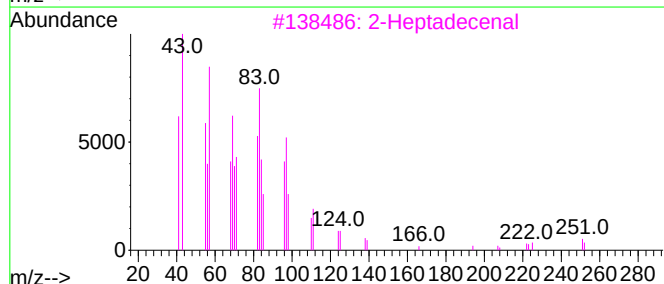
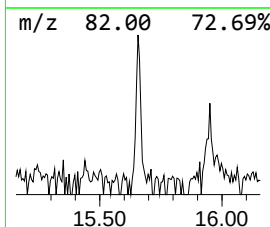
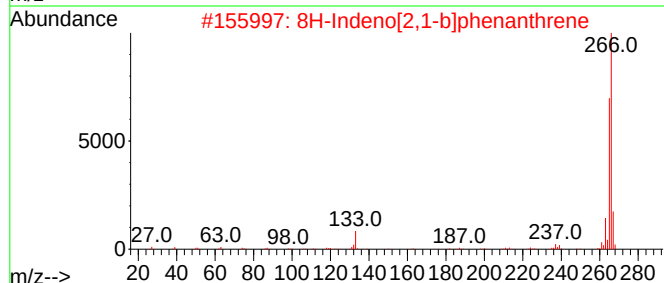
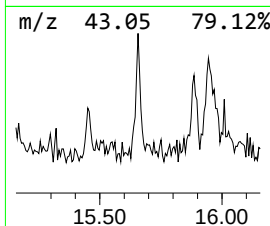
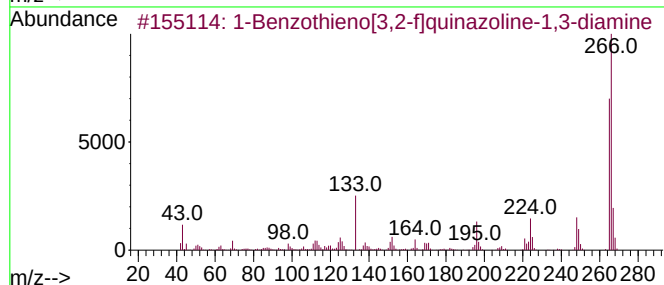
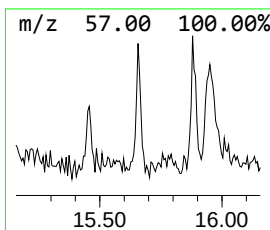
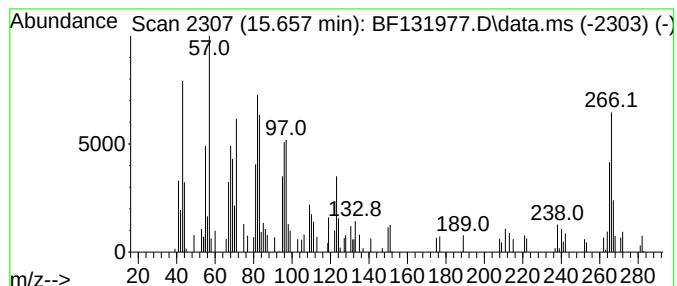
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown15.656 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.656	2.91 ng	61260	Perylene-d12	15.439

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Benzothieno[3,2-f]quinazoline-...	266	C14H10N4S	065642-92-4	25
2		8H-Indeno[2,1-b]phenanthrene	266	C21H14	000241-28-1	22
3		2-Heptadecenal	252	C17H32O	1000143-48-6	18
4		16-Heptadecenal	252	C17H32O	1010144-57-9	15
5		14-Heptadecenal	252	C17H32O	1010144-58-0	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
Operator : CG\JU
Sample : 01065-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-210

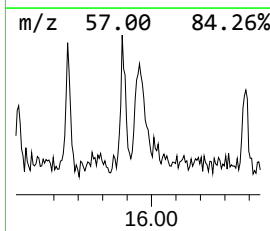
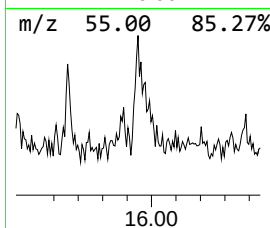
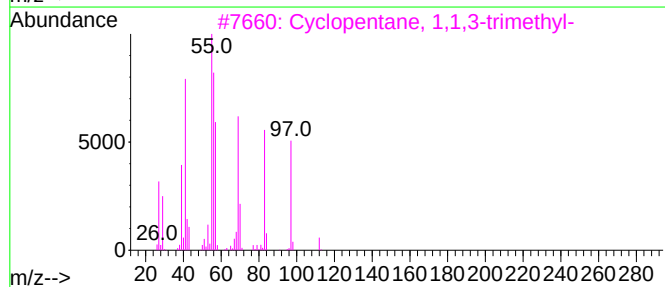
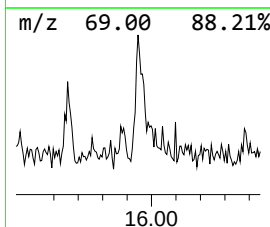
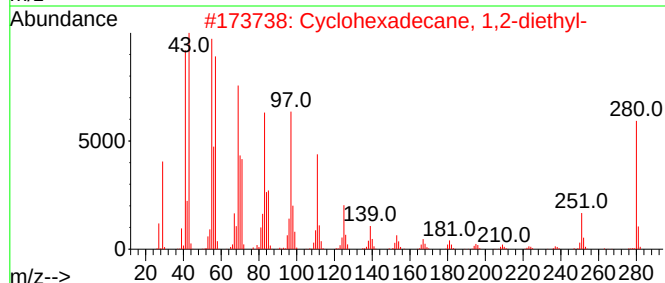
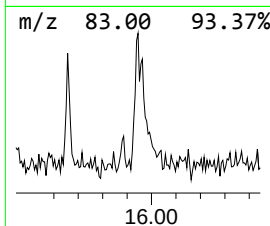
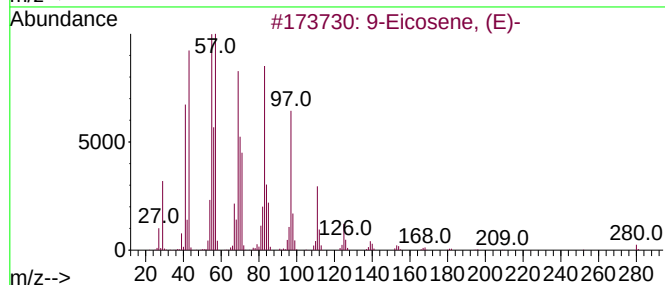
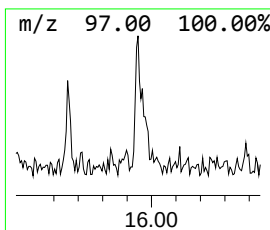
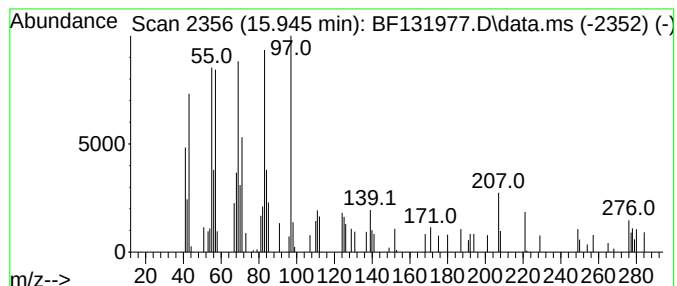
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9-Eicosene, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	2.78 ng	58545	Perylene-d12	15.439

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Eicosene, (E)-	280	C20H40	074685-29-3	74
2		Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	74
3		Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	68
4		Cyclododecane	168	C12H24	000294-62-2	58
5		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	000547-60-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131977.D
Acq On : 09 Jan 2023 22:39
Operator : CG\JU
Sample : 01065-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VNJ-210

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.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
Propane, 1,1-di...	2.128	2.7	ng	57397	1	6.781	428791	20.0
2-Pentanone, 4-...	4.987	37.6	ng	805168	1	6.781	428791	20.0
Benzophenone	10.551	4.1	ng	130871	3	9.833	639378	20.0
Phenanthrene, 1...	11.857	2.7	ng	95122	4	11.327	708944	20.0
4H-Cyclopenta[d...	11.939	3.1	ng	111280	4	11.327	708944	20.0
11H-Benzo[b]flu...	12.998	2.0	ng	76624	5	13.974	760406	20.0
Pyrene, 2-methyl-	13.104	2.3	ng	87748	5	13.974	760406	20.0
Heptadecyl hept...	13.816	6.5	ng	245994	5	13.974	760406	20.0
Perylene	15.121	3.0	ng	63346	6	15.439	421074	20.0
Benzo[e]pyrene	15.315	7.3	ng	153959	6	15.439	421074	20.0
unknown15.656	15.656	2.9	ng	61260	6	15.439	421074	20.0
9-Eicosene, (E)-	15.945	2.8	ng	58545	6	15.439	421074	20.0