

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
 Data File : BF129929.D
 Acq On : 20 Aug 2022 01:51
 Operator : CG\JU
 Sample : N4217-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP2-0815

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-BF081422.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129929.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.004	460	462	475	rBV	107168	154757	4.82%	0.594%
2	5.422	529	533	556	rBV	2894926	2767359	86.15%	10.628%
3	6.440	702	706	721	rBV	2922978	2827011	88.01%	10.857%
4	6.787	761	765	774	rBV	949226	832439	25.92%	3.197%
5	7.357	857	862	871	rBV	1773264	1763077	54.89%	6.771%
6	8.069	979	983	986	rBV	1340756	1152836	35.89%	4.428%
7	9.139	1161	1165	1168	rBV	3696275	3212099	100.00%	12.336%
8	9.822	1277	1281	1284	rBV	1758900	1353922	42.15%	5.200%
9	10.616	1412	1416	1429	rBV	2600643	2231372	69.47%	8.570%
10	11.310	1531	1534	1537	rBV	1596814	1363654	42.45%	5.237%
11	11.333	1537	1538	1543	rVB	209085	143572	4.47%	0.551%
12	11.686	1596	1598	1601	rVB	49239	40756	1.27%	0.157%
13	11.816	1617	1620	1621	rBV	110406	101248	3.15%	0.389%
14	11.839	1621	1624	1628	rVB2	132533	162928	5.07%	0.626%
15	11.922	1635	1638	1641	rBV2	123253	132021	4.11%	0.507%
16	11.945	1641	1642	1647	rVB	87622	73529	2.29%	0.282%
17	12.110	1666	1670	1674	rBV2	43645	51345	1.60%	0.197%
18	12.363	1709	1713	1715	rBV	105296	100104	3.12%	0.384%
19	12.398	1715	1719	1721	rVV3	84430	107831	3.36%	0.414%
20	12.463	1725	1730	1736	rVB3	45326	87285	2.72%	0.335%
21	12.527	1737	1741	1745	rBV	299178	255582	7.96%	0.982%
22	12.757	1777	1780	1786	rVB	402278	358226	11.15%	1.376%
23	12.810	1786	1789	1792	rVB	42870	39227	1.22%	0.151%
24	12.898	1799	1804	1807	rBV	3108000	2946630	91.74%	11.317%
25	12.975	1813	1817	1823	rBV5	50052	88519	2.76%	0.340%
26	13.063	1829	1832	1834	rBV3	39523	50884	1.58%	0.195%
27	13.086	1834	1836	1838	rVB2	51475	43585	1.36%	0.167%
28	13.192	1850	1854	1858	rVB	77957	72388	2.25%	0.278%
29	13.286	1866	1870	1872	rVB	57581	50332	1.57%	0.193%
30	13.316	1872	1875	1878	rBV	60903	58611	1.82%	0.225%
31	13.604	1921	1924	1927	rVB	36199	37641	1.17%	0.145%
32	13.710	1937	1942	1944	rBV2	46838	56628	1.76%	0.217%
33	13.786	1951	1955	1969	rVB5	186312	467713	14.56%	1.796%
34	13.957	1979	1984	1987	rBV	1001759	997670	31.06%	3.832%
35	13.986	1987	1989	1992	rVV	131044	117338	3.65%	0.451%
36	14.045	1996	1999	2005	rVV	74910	78470	2.44%	0.301%

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
 Data File : BF129929.D
 Acq On : 20 Aug 2022 01:51
 Operator : CG\JU
 Sample : N4217-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP2-0815

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-BF081422.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.474	2070	2072	2079	rVB2	43486	59093	1.84%	0.227%
38	14.698	2107	2110	2113	rVB	39233	40081	1.25%	0.154%
39	14.998	2157	2161	2163	rBV	67102	97344	3.03%	0.374%
40	15.021	2163	2165	2173	rVB2	92910	109806	3.42%	0.422%
41	15.298	2208	2212	2217	rVB	55743	70785	2.20%	0.272%
42	15.357	2218	2222	2225	rBV	75863	92769	2.89%	0.356%
43	15.416	2228	2232	2240	rVB	643736	747294	23.26%	2.870%
44	15.810	2295	2299	2306	rVB	57656	79317	2.47%	0.305%
45	16.292	2377	2381	2391	rVB3	56894	110637	3.44%	0.425%
46	16.874	2473	2480	2489	rVB3	56931	147359	4.59%	0.566%
47	17.327	2553	2557	2565	rVB3	26297	49058	1.53%	0.188%
48	17.533	2588	2592	2599	rVB3	26888	55552	1.73%	0.213%

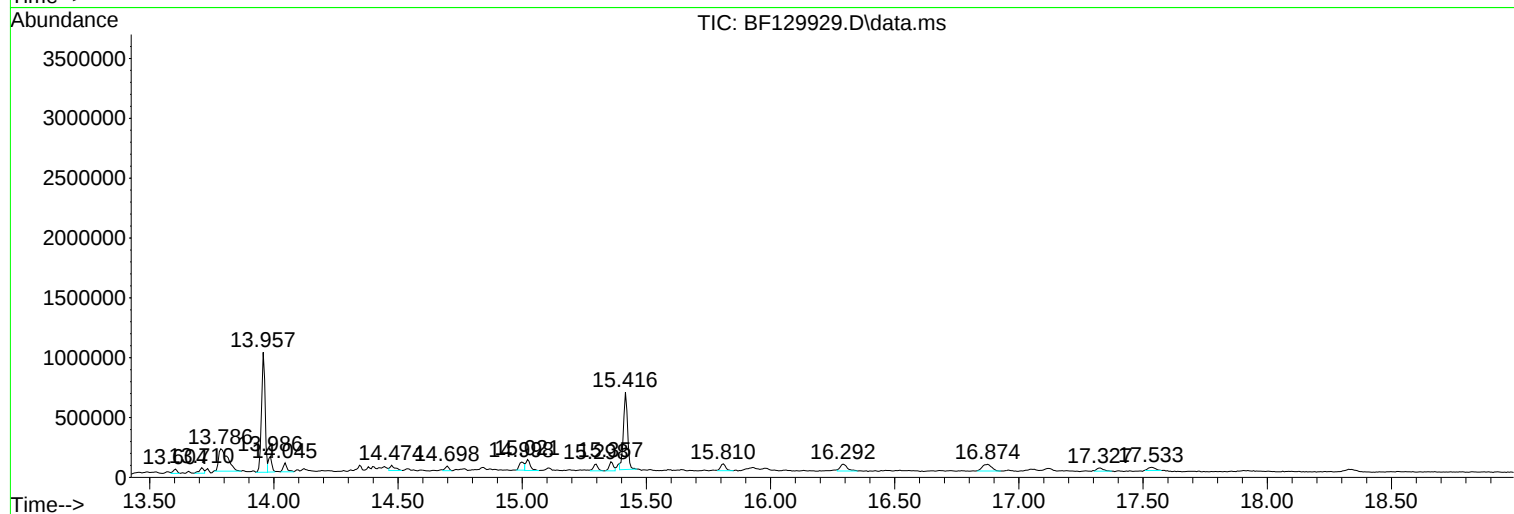
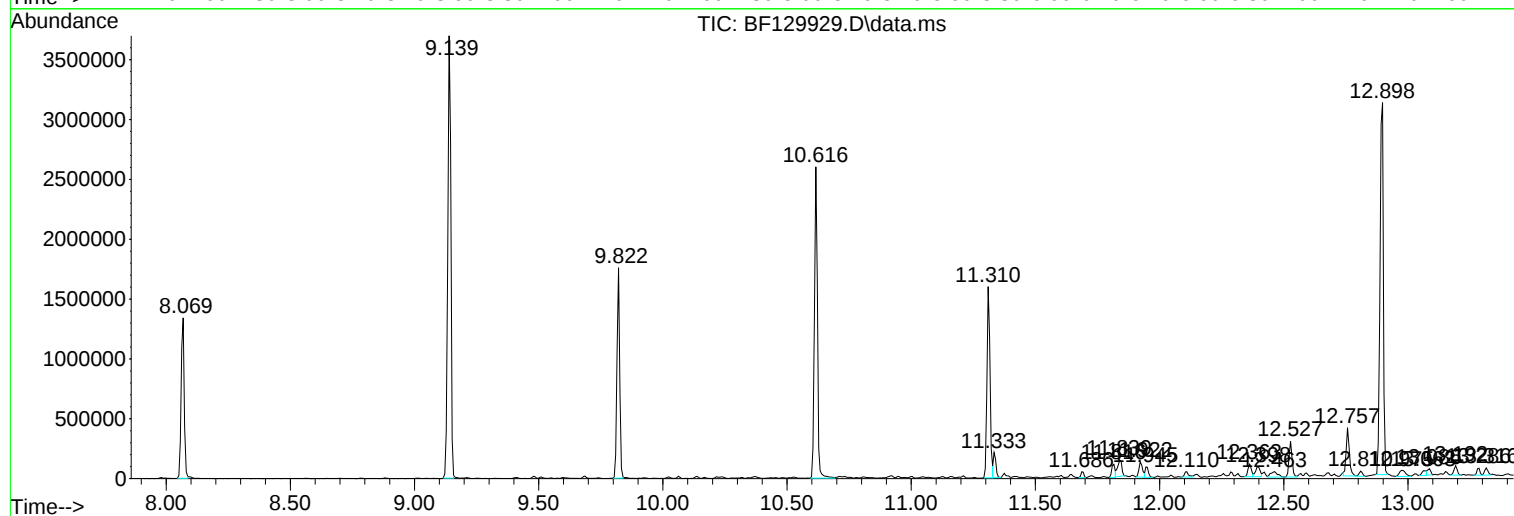
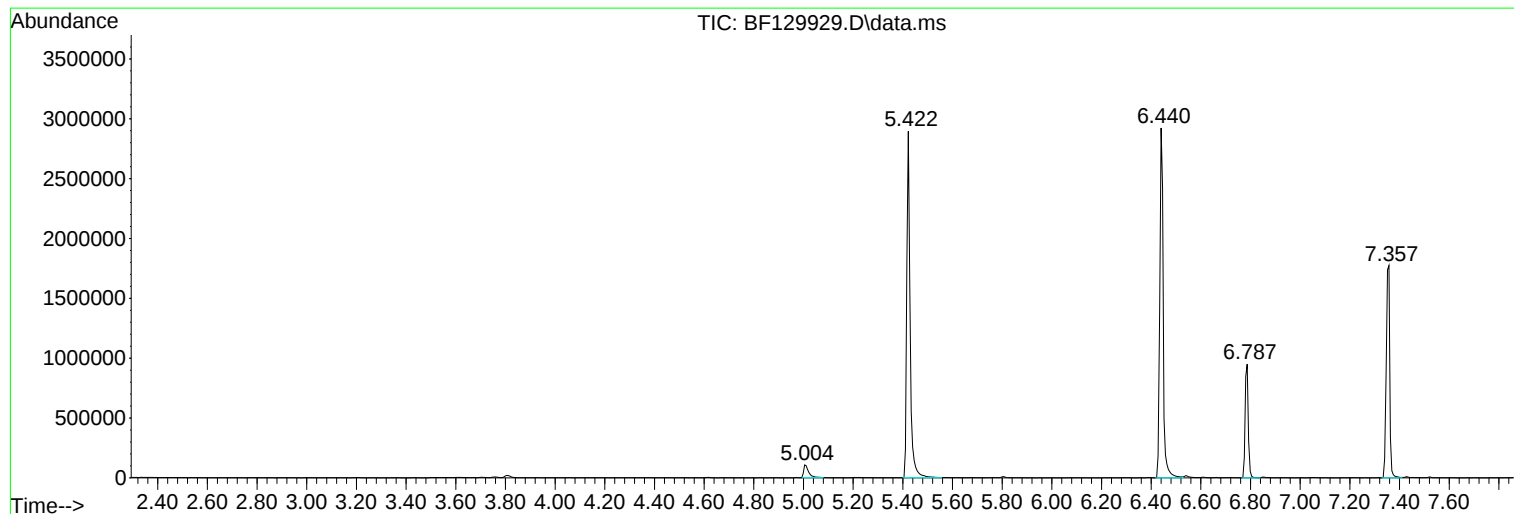
Sum of corrected areas: 26037684

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP2-0815

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP2-0815

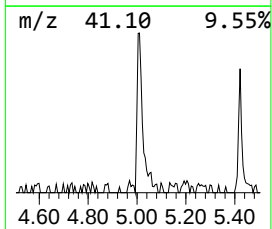
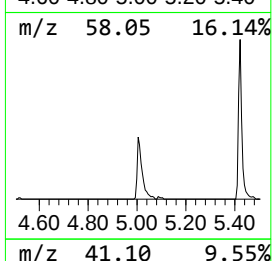
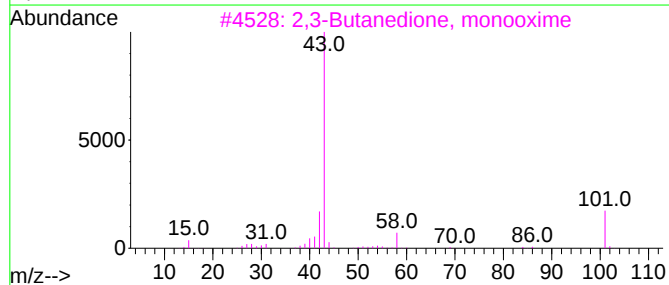
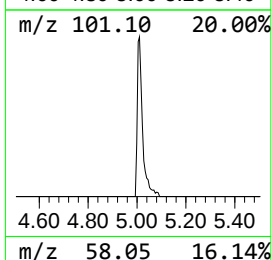
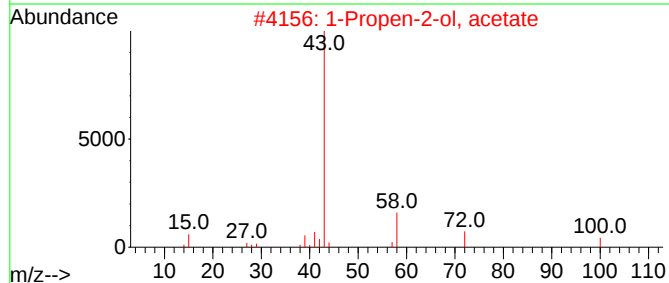
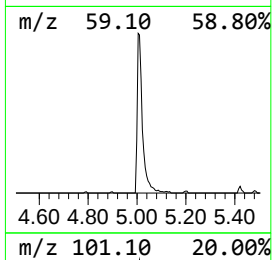
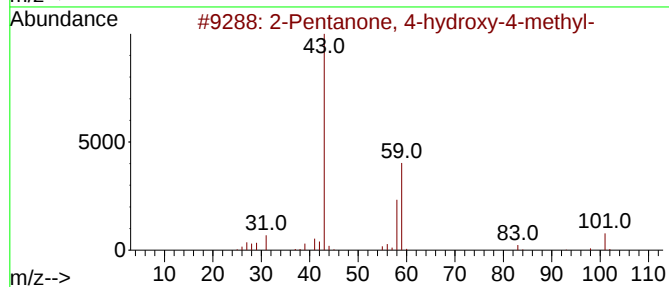
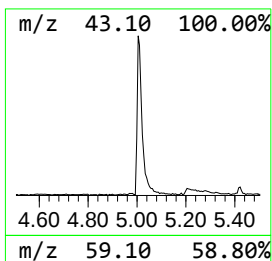
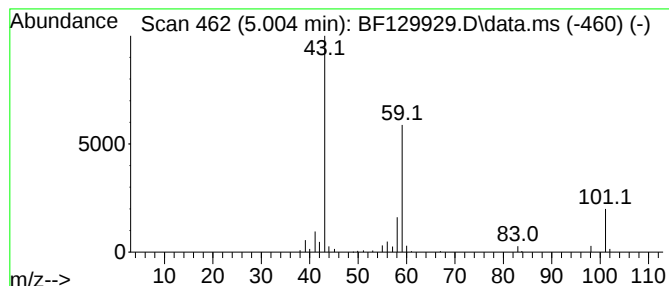
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.004	3.72 ng	154757	1,4-Dichlorobenzene-d4	6.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42		
2	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP2-0815

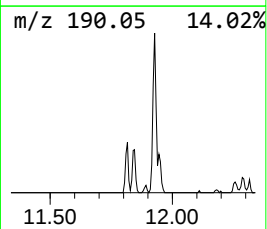
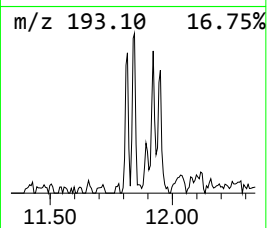
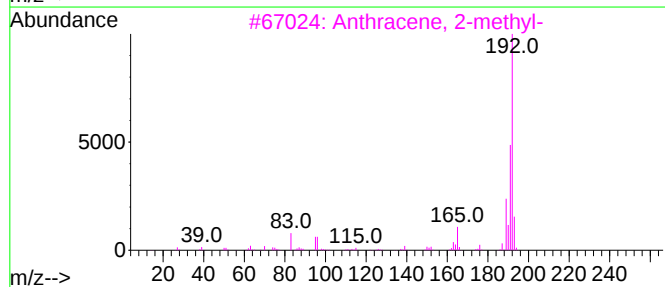
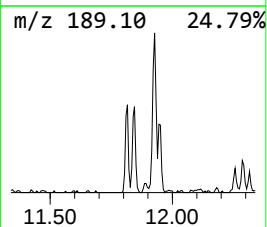
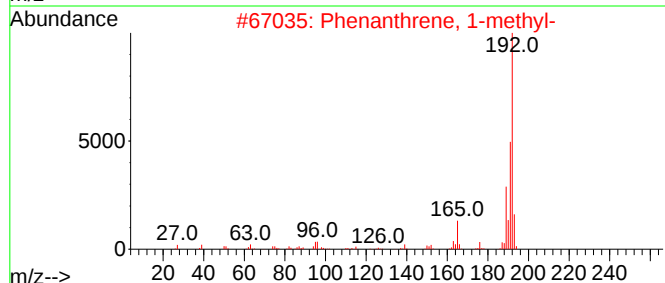
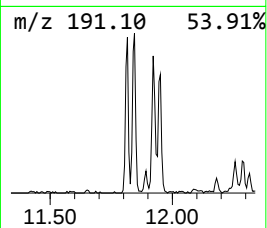
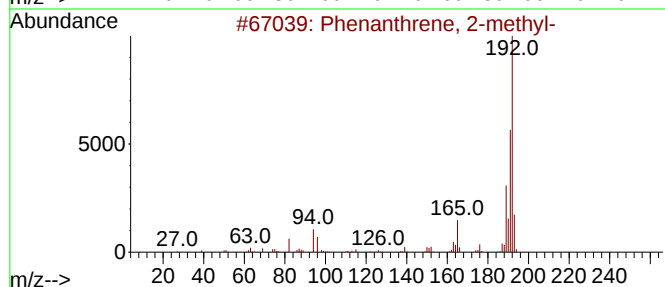
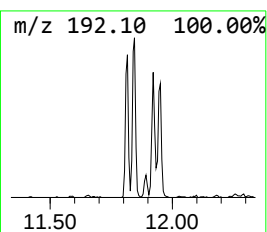
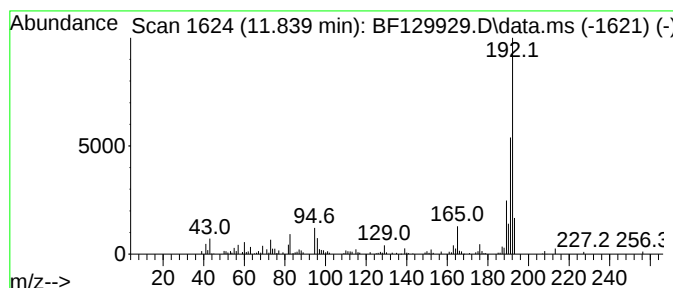
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.839	2.39 ng	162928	Phenanthrene-d10	11.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP2-0815

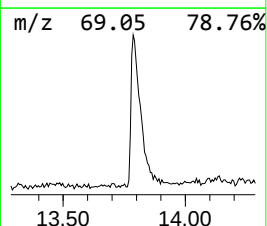
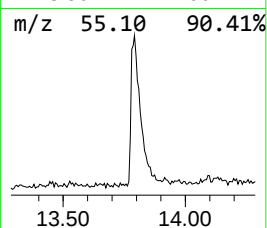
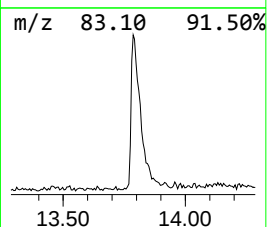
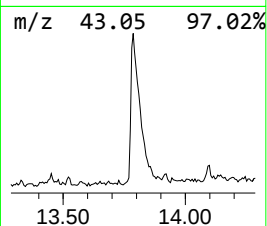
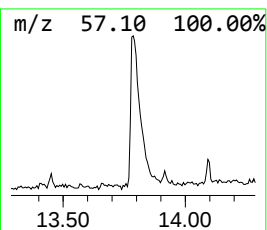
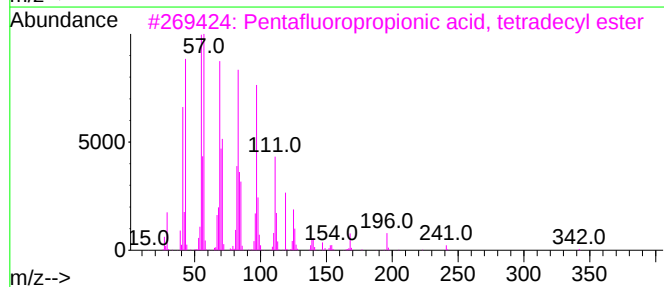
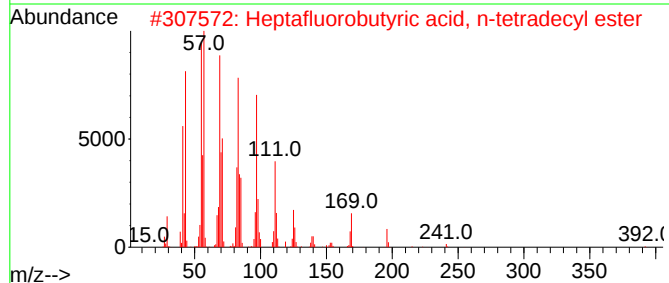
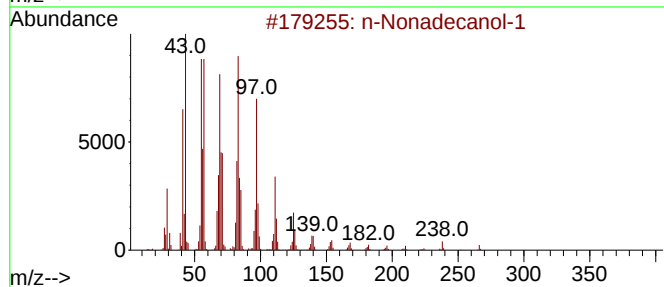
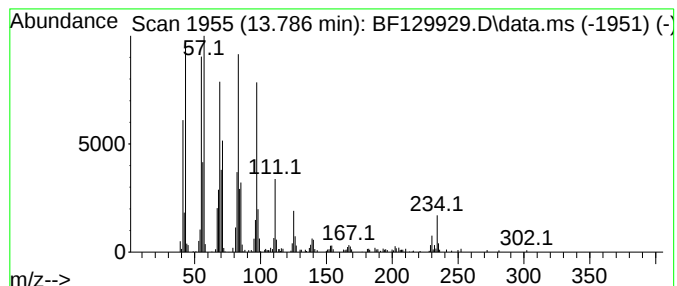
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Nonadecanol-1 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.786	9.38 ng	467713	Chrysene-d12	13.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
3			Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	94
4			1-Heneicosanol	312	C21H44O	015594-90-8	93
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP2-0815

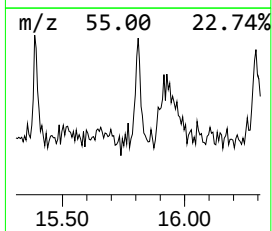
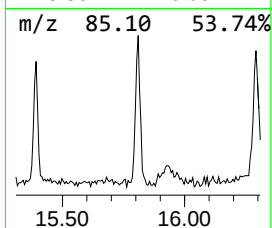
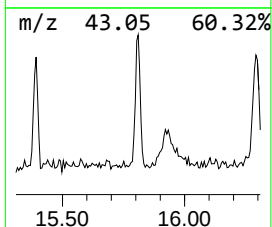
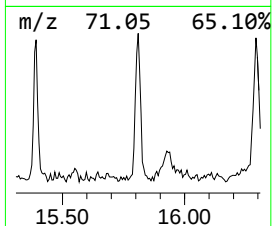
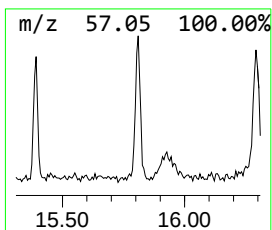
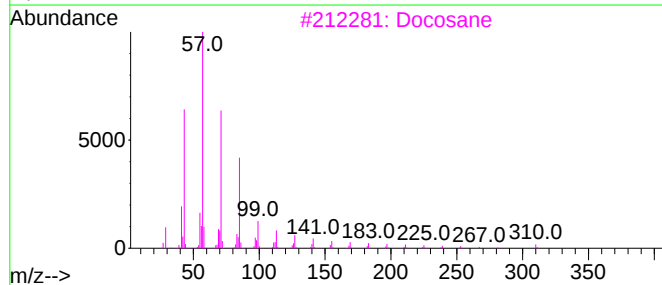
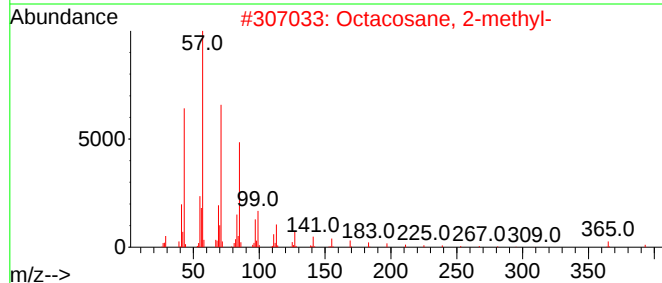
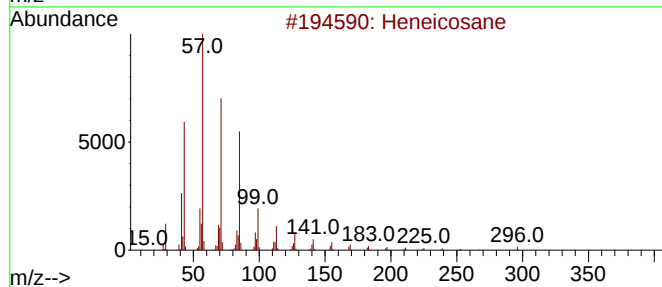
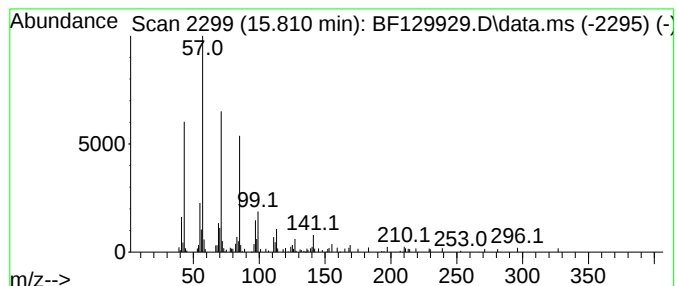
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.810	2.12 ng	79317	Perylene-d12	15.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	98
2			Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
3			Docosane	310	C22H46	000629-97-0	86
4			Heptadecane	240	C17H36	000629-78-7	86
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP2-0815

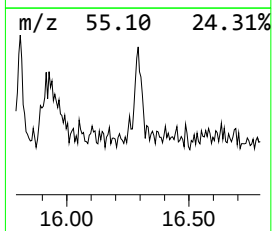
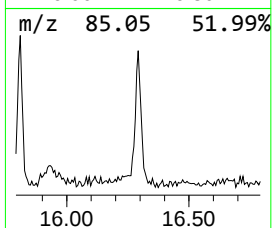
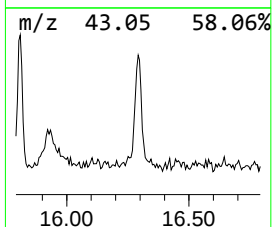
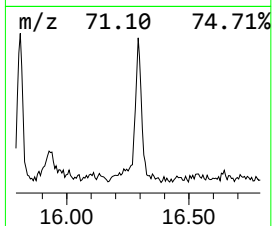
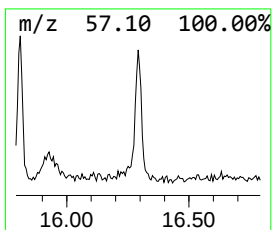
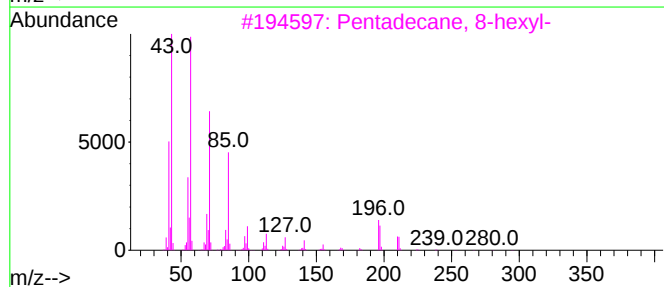
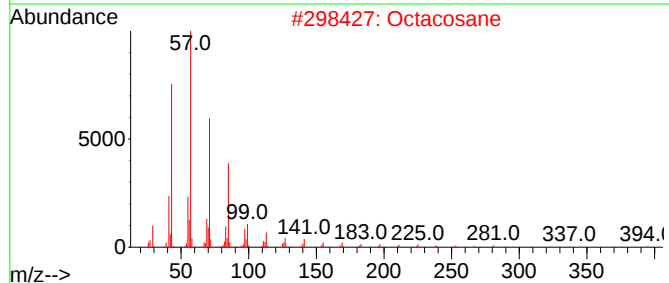
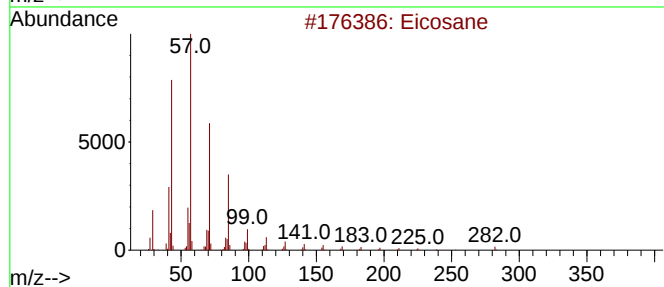
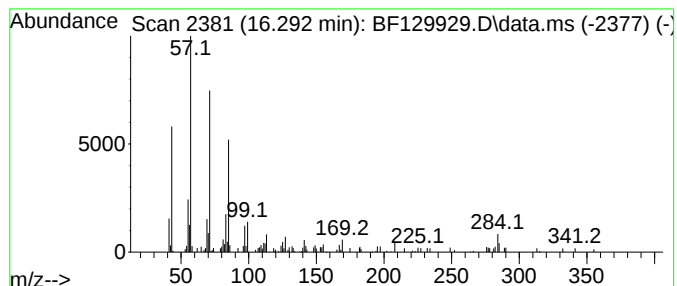
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Eicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.292	2.96 ng	110637	Perylene-d12	15.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			Octacosane	394	C28H58	000630-02-4	87
3			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	86
4			Hexacosane	366	C26H54	000630-01-3	86
5			Docosane, 11-butyl-	366	C26H54	013475-76-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

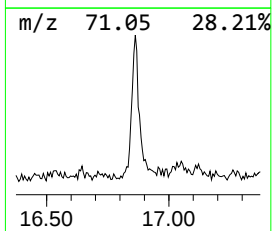
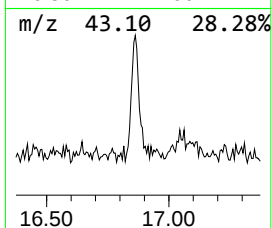
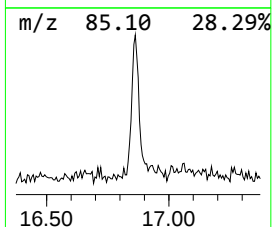
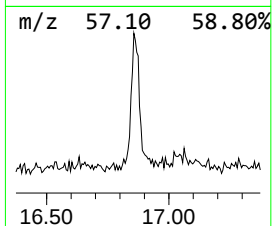
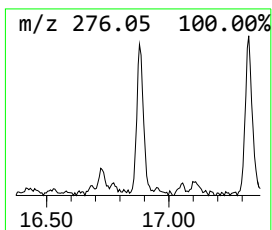
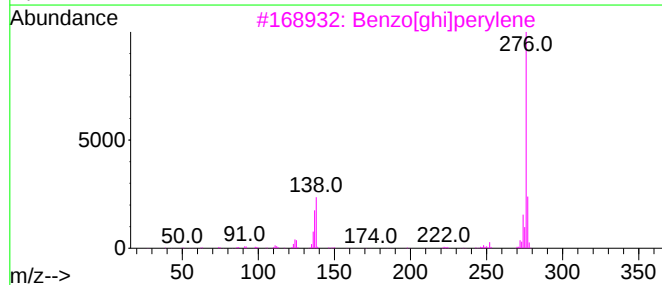
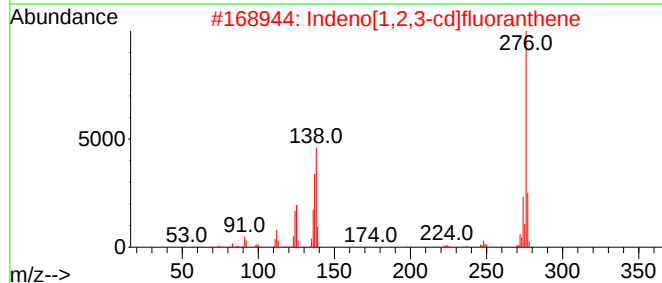
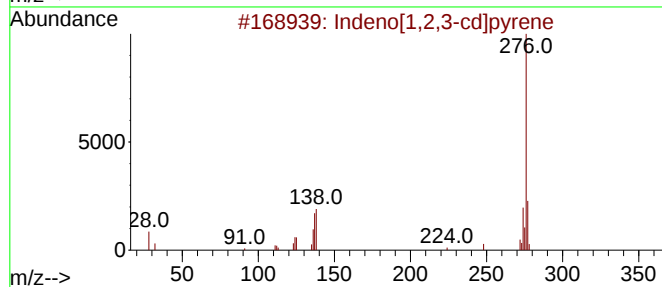
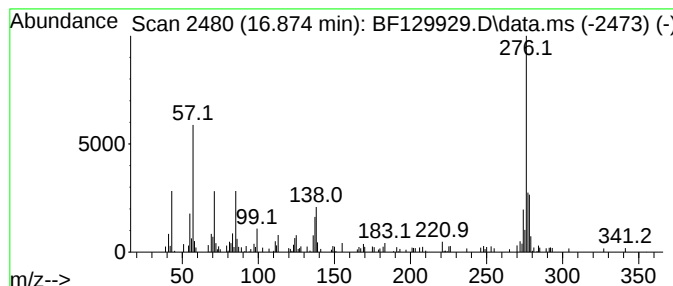
Instrument :
BNA_F
ClientSampleId :
TP2-0815

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Indeno[1,2,3-cd]fluoranthene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.874	3.94 ng	147359	Perylene-d12	15.416	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	93
2	Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	93
3	Benzo[ghi]perylene	276	C22H12	000191-24-2	70
4	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	64
5	Propionamide, N-[1-(2-chloropyri...	349	C11H10ClF6N3O	1000269-44-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081922\
Data File : BF129929.D
Acq On : 20 Aug 2022 01:51
Operator : CG\JU
Sample : N4217-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP2-0815

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.004	3.7	ng	154757	1	6.787	832439	20.0
Phenanthrene, 2...	11.839	2.4	ng	162928	4	11.310	1363650	20.0
n-Nonadecanol-1	13.786	9.4	ng	467713	5	13.957	997670	20.0
Heneicosane	15.810	2.1	ng	79317	6	15.416	747294	20.0
Eicosane	16.292	3.0	ng	110637	6	15.416	747294	20.0
Indeno[1,2,3-cd...	16.874	3.9	ng	147359	6	15.416	747294	20.0