

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
 Data File : BF137724.D
 Acq On : 24 Apr 2024 18:18
 Operator : CG\JU
 Sample : P2245-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 YARD-3

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-BF042324.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF137724.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.222	16	23	30	rVB3	77259	175523	3.05%	0.379%
2	2.440	54	60	69	rVB	394878	515479	8.95%	1.114%
3	3.698	272	274	287	rBV	91987	147407	2.56%	0.318%
4	5.681	607	611	615	rVB	4603984	4666553	81.00%	10.081%
5	6.669	773	779	782	rBV	4903424	4807127	83.44%	10.385%
6	7.063	842	846	849	rBV	1721259	1487842	25.83%	3.214%
7	7.622	936	941	944	rBV	3292994	3036246	52.70%	6.559%
8	7.857	978	981	985	rVB	106186	86768	1.51%	0.187%
9	8.228	1041	1044	1049	rBV	226670	177182	3.08%	0.383%
10	8.345	1060	1064	1067	rBV	2487446	2006275	34.83%	4.334%
11	8.645	1111	1115	1118	rVB	122525	110903	1.93%	0.240%
12	9.416	1242	1246	1250	rBV	6258863	5760865	100.00%	12.445%
13	10.104	1359	1363	1366	rBV	2770201	2211598	38.39%	4.778%
14	10.139	1366	1369	1371	rVB	132552	114476	1.99%	0.247%
15	10.186	1371	1377	1384	rBV4	120691	179341	3.11%	0.387%
16	10.310	1394	1398	1401	rVB	123459	112649	1.96%	0.243%
17	10.657	1453	1457	1462	rBV	200455	197618	3.43%	0.427%
18	10.898	1493	1498	1501	rBV	3845308	3498251	60.72%	7.557%
19	11.204	1542	1550	1552	rBV3	50581	64790	1.12%	0.140%
20	11.457	1587	1593	1597	rBV4	34627	80114	1.39%	0.173%
21	11.498	1597	1600	1604	rVB	99851	83823	1.46%	0.181%
22	11.598	1614	1617	1620	rBV	2379952	2348272	40.76%	5.073%
23	11.621	1620	1621	1625	rVV	1500654	1147671	19.92%	2.479%
24	11.674	1625	1630	1633	rVV	606703	521124	9.05%	1.126%
25	11.704	1633	1635	1638	rVB	91534	81281	1.41%	0.176%
26	11.827	1648	1656	1661	rBV	317923	344777	5.98%	0.745%
27	12.104	1699	1703	1705	rBV	363881	326344	5.66%	0.705%
28	12.127	1705	1707	1711	rVV2	217599	199826	3.47%	0.432%
29	12.174	1711	1715	1718	rVB2	80733	85605	1.49%	0.185%
30	12.216	1718	1722	1724	rBV2	265038	276957	4.81%	0.598%
31	12.239	1724	1726	1728	rVB	99681	78164	1.36%	0.169%
32	12.363	1744	1747	1749	rBV2	70100	64538	1.12%	0.139%
33	12.398	1749	1753	1755	rVV2	89043	80808	1.40%	0.175%
34	12.651	1793	1796	1799	rBV	63619	63237	1.10%	0.137%
35	12.815	1820	1824	1828	rBV	1255940	1089182	18.91%	2.353%
36	12.892	1833	1837	1842	rVB2	58845	86002	1.49%	0.186%

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
YARD-3

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-BF042324.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.051	1861	1864	1869	rVB	952818	875355	15.19%	1.891%
38	13.186	1883	1887	1890	rBV	5723733	4850695	84.20%	10.479%
39	13.257	1895	1899	1904	rBV2	49548	86324	1.50%	0.186%
40	13.374	1916	1919	1922	rVB	183833	159746	2.77%	0.345%
41	13.445	1927	1931	1933	rBV	162853	155382	2.70%	0.336%
42	13.480	1933	1937	1940	rVV2	61351	69565	1.21%	0.150%
43	13.998	2020	2025	2027	rBV	60446	59953	1.04%	0.130%
44	14.068	2032	2037	2042	rVV5	118228	157631	2.74%	0.341%
45	14.251	2063	2068	2070	rBV2	1519195	1565194	27.17%	3.381%
46	14.274	2070	2072	2077	rVB	249277	223212	3.87%	0.482%
47	15.345	2250	2254	2258	rBV	105977	185379	3.22%	0.400%
48	15.680	2307	2311	2318	rVB	64232	92343	1.60%	0.199%
49	15.751	2318	2323	2329	rBV	89912	116179	2.02%	0.251%
50	15.821	2329	2335	2348	rBV	1003229	1377590	23.91%	2.976%

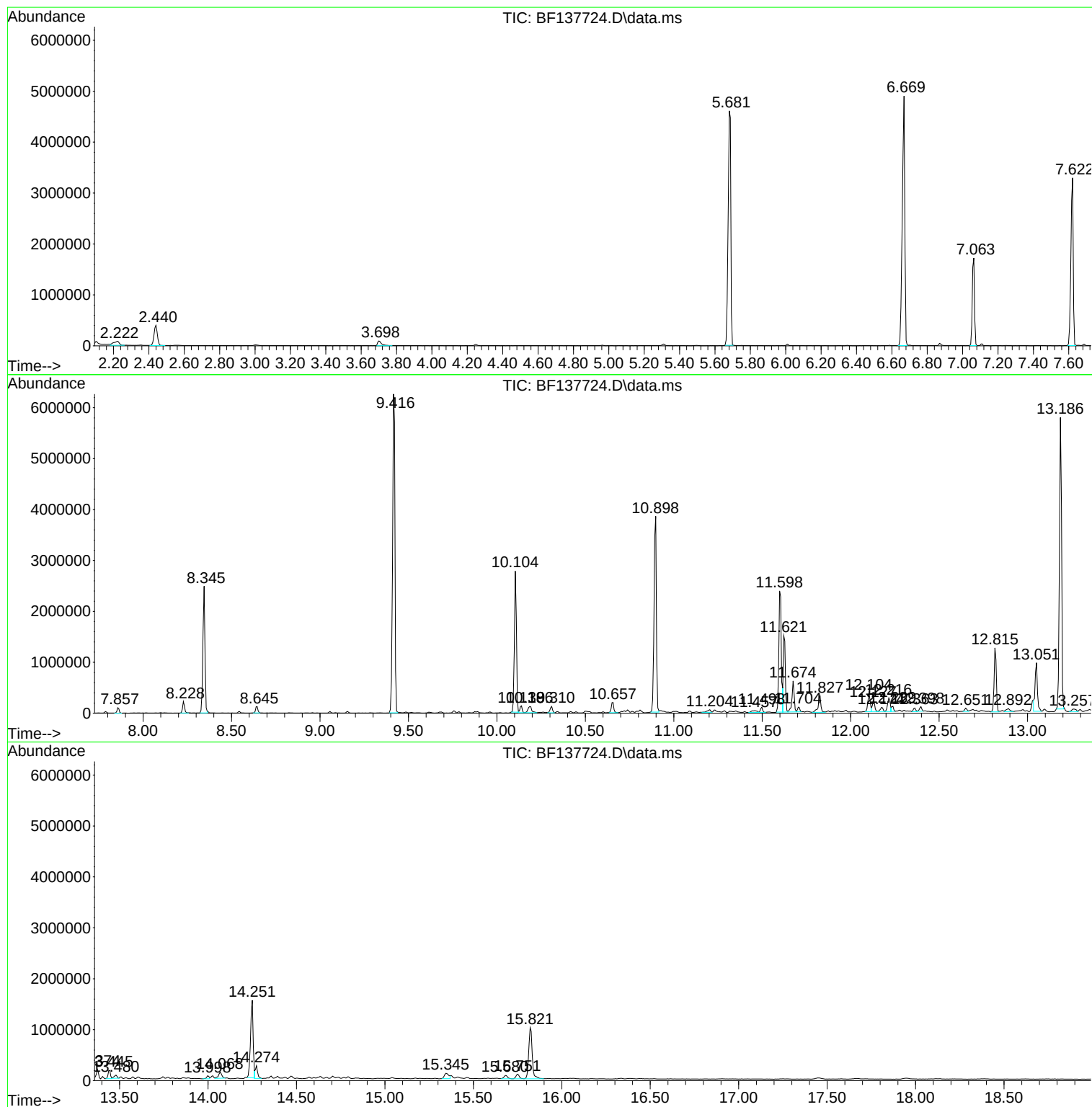
Sum of corrected areas: 46289166

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
YARD-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.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\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
YARD-3

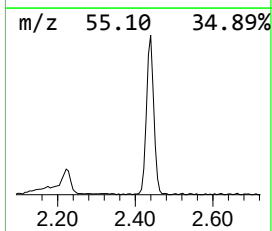
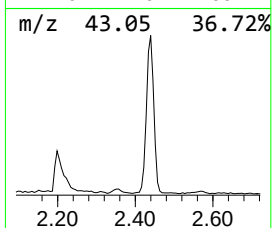
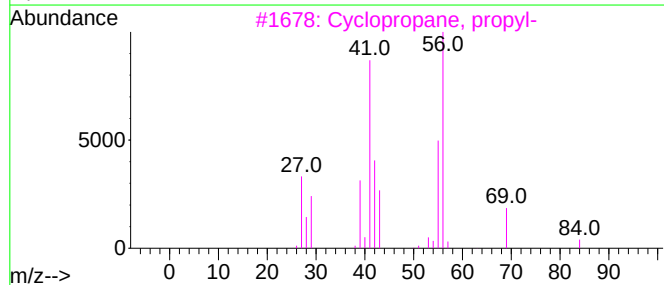
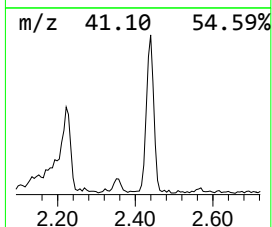
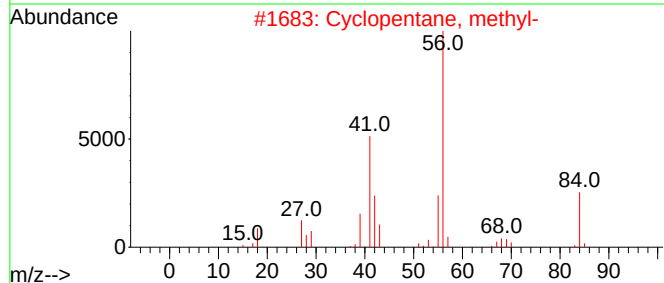
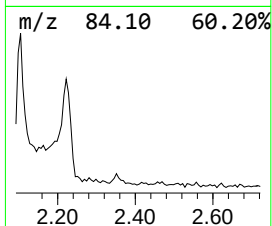
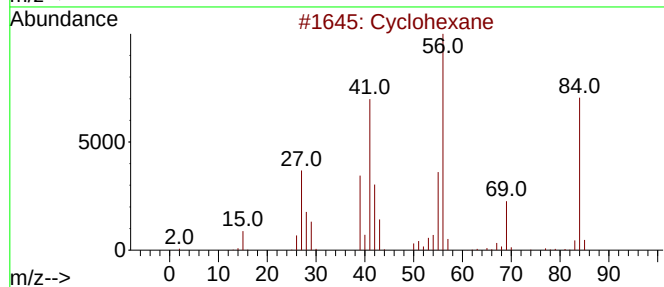
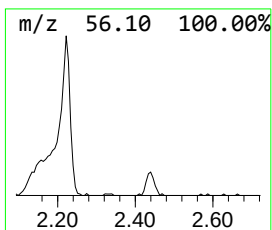
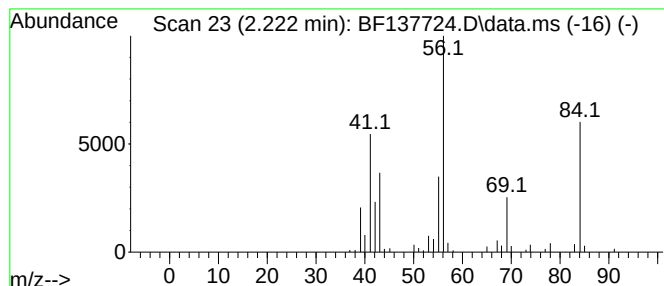
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.222	2.36 ng	175523	1,4-Dichlorobenzene-d4	7.063

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane	84	C6H12	000110-82-7	90
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	56
3		Cyclopropane, propyl-	84	C6H12	002415-72-7	50
4		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	45
5		Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
YARD-3

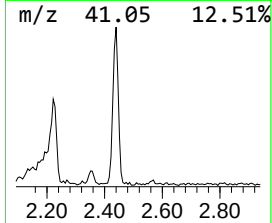
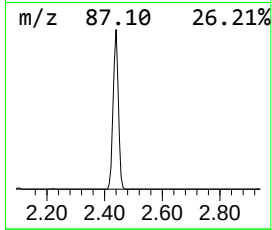
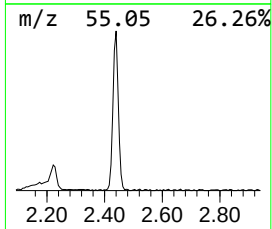
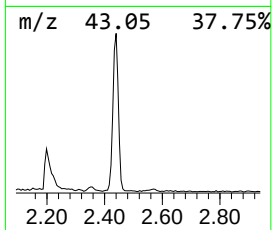
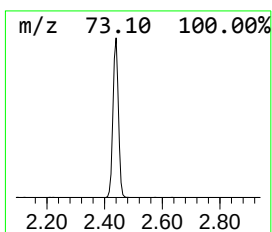
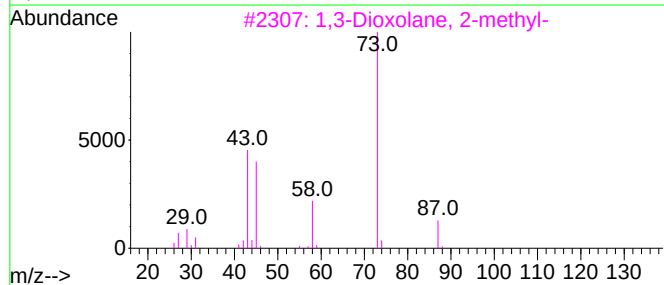
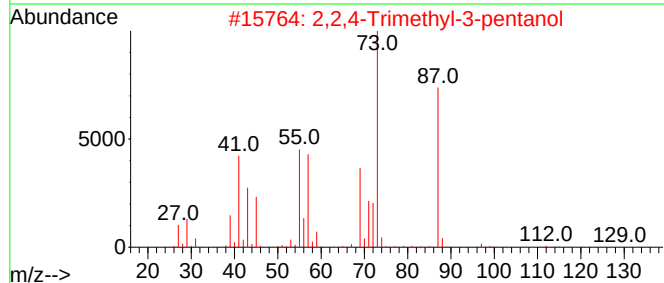
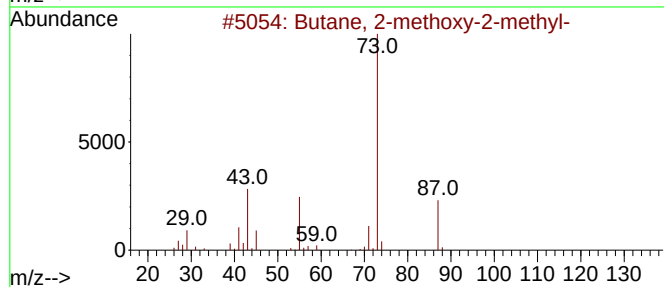
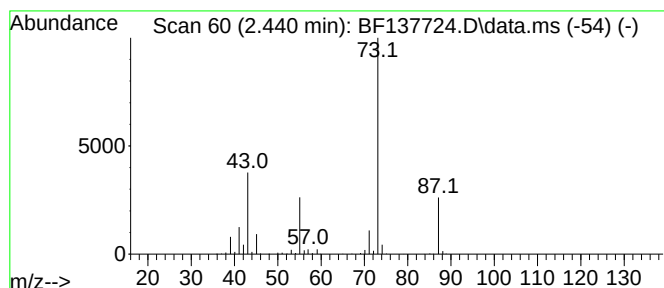
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.440	6.93 ng	515479	1,4-Dichlorobenzene-d4	7.063

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5		1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

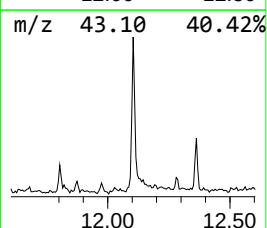
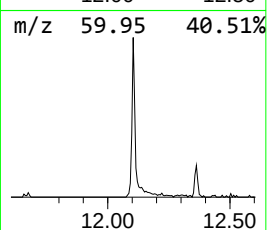
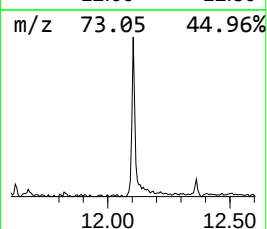
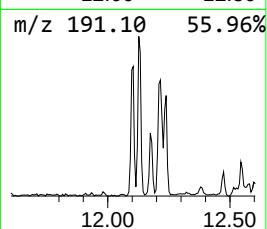
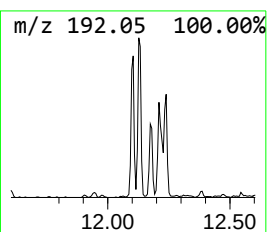
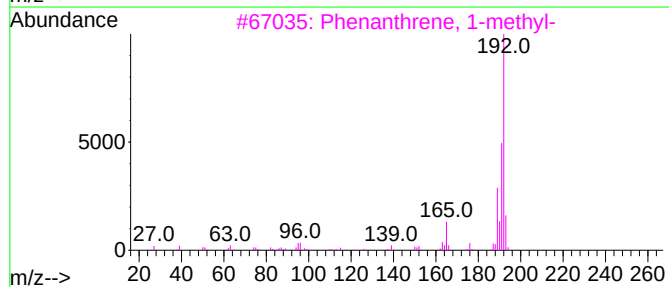
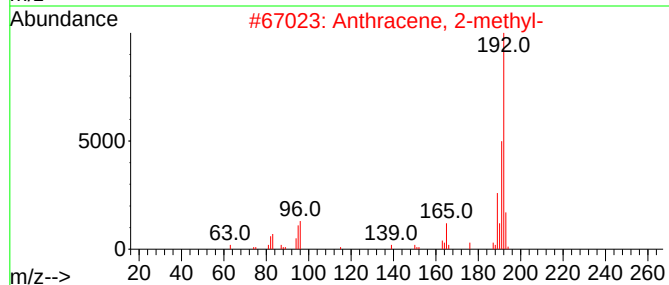
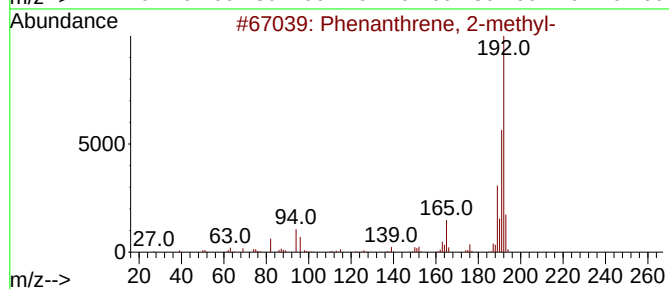
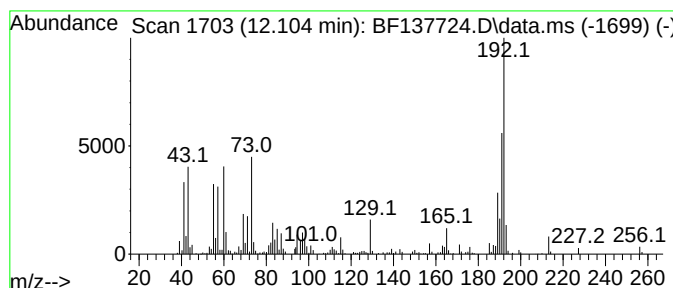
Instrument :
BNA_F
ClientSampleId :
YARD-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.104	2.78 ng	326344	Phenanthrene-d10	11.598	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	89
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	83
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	83
5	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

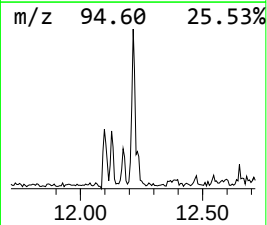
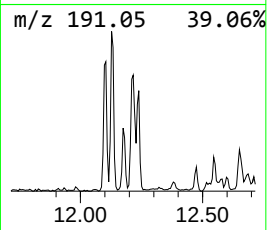
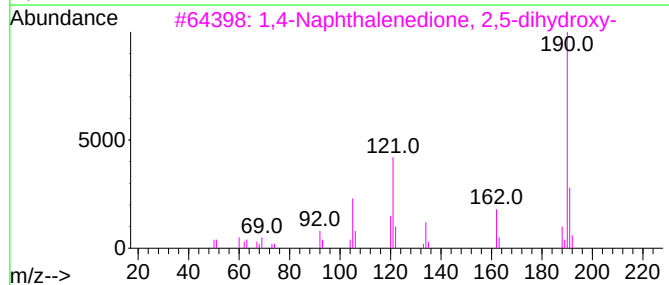
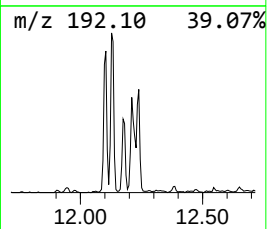
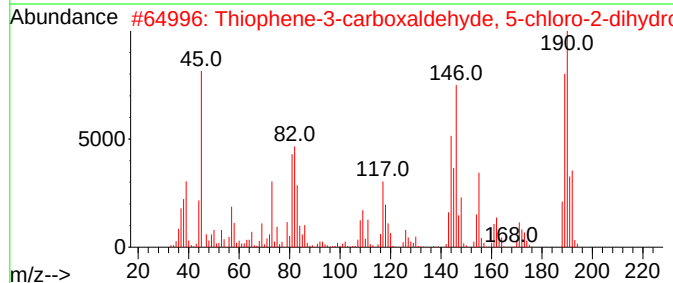
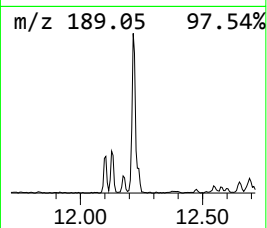
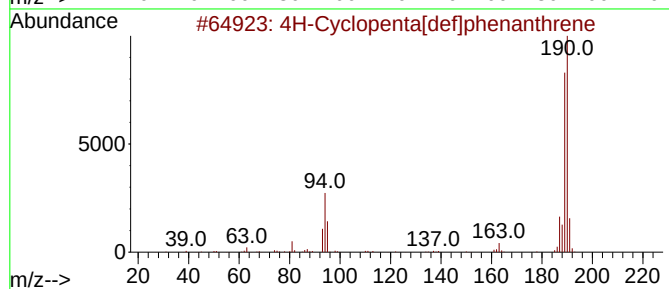
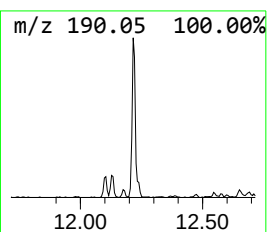
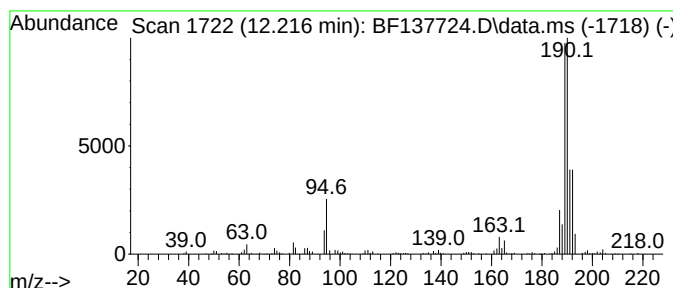
Instrument :
BNA_F
ClientSampleId :
YARD-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.216	2.36 ng	276957	Phenanthrene-d10	11.598	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
3	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	25
5	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
YARD-3

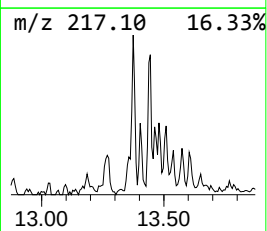
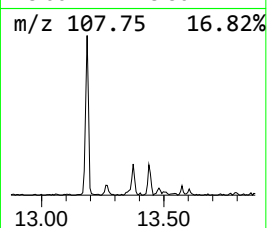
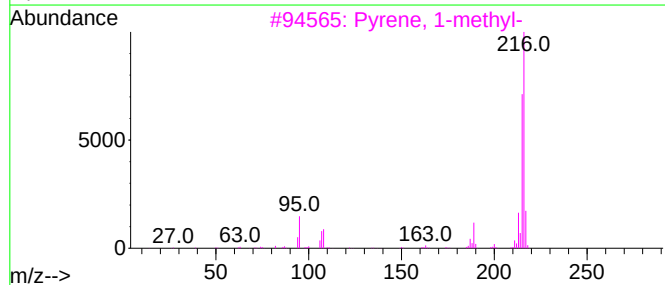
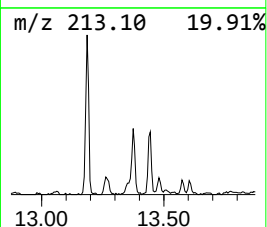
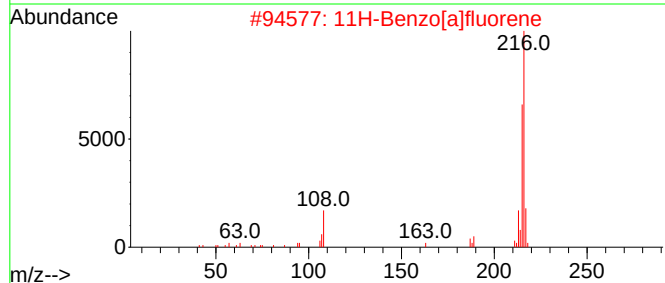
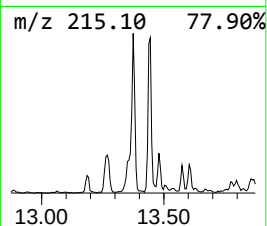
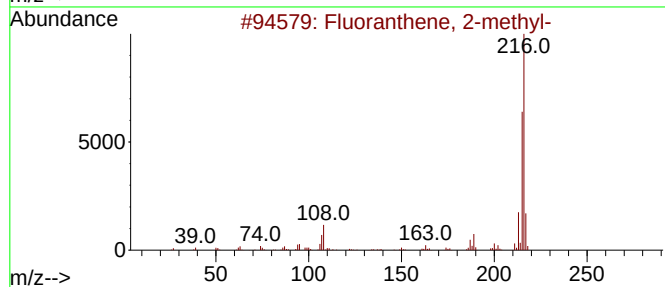
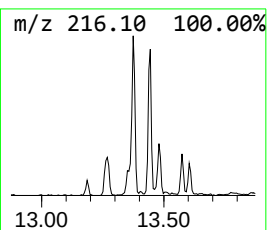
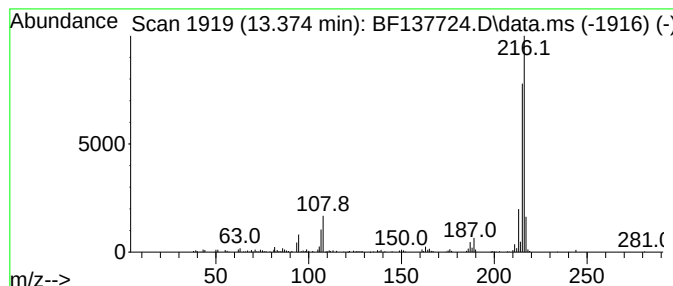
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.374	2.04 ng	159746	Chrysene-d12	14.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
5			Pyrene, 4-methyl-	216	C17H12	003353-12-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
YARD-3

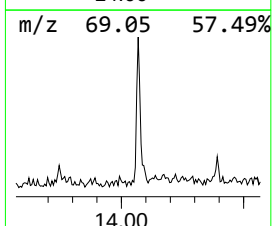
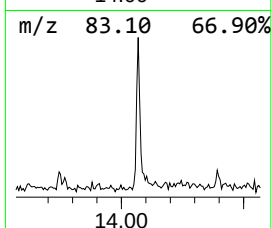
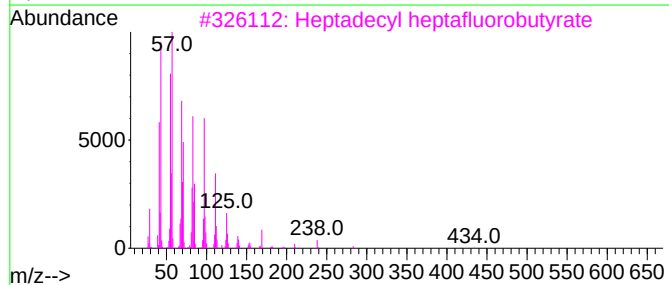
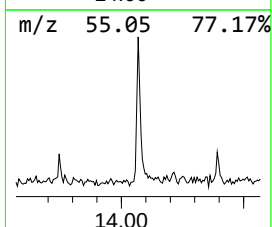
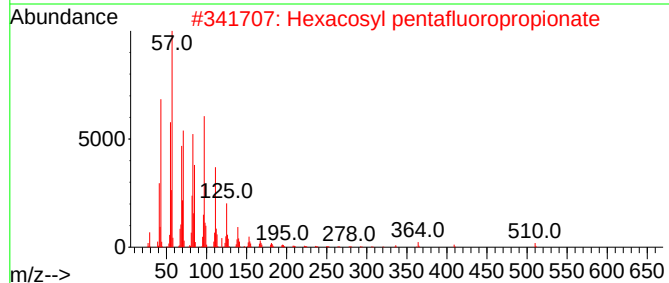
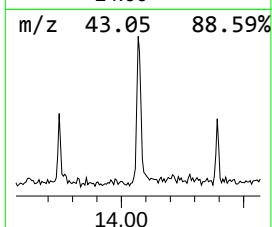
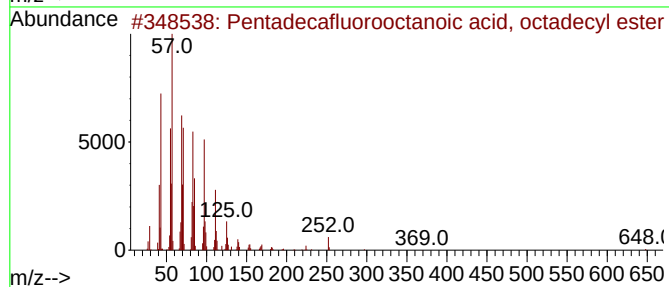
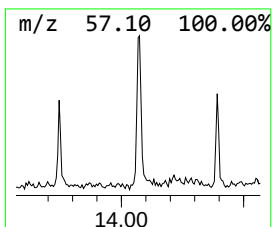
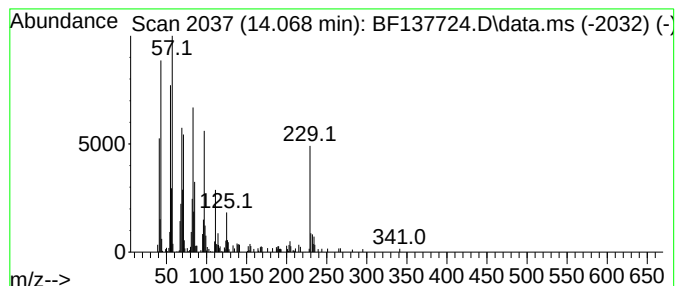
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecafluorooctanoic aci... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.068	2.01 ng	157631	Chrysene-d12	14.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	86
2			Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	70
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	70
4			Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	70
5			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042424\
Data File : BF137724.D
Acq On : 24 Apr 2024 18:18
Operator : CG\JU
Sample : P2245-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
YARD-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042324.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
Cyclohexane	2.222	2.4	ng	175523	1	7.063	1487840	20.0
Butane, 2-metho...	2.440	6.9	ng	515479	1	7.063	1487840	20.0
Phenanthrene, 2...	12.104	2.8	ng	326344	4	11.598	2348270	20.0
4H-Cyclopenta[d...	12.216	2.4	ng	276957	4	11.598	2348270	20.0
Fluoranthene, 2...	13.374	2.0	ng	159746	5	14.251	1565190	20.0
Pentadecafluoro...	14.068	2.0	ng	157631	5	14.251	1565190	20.0