

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090319\
 Data File : BG042693.D
 Acq On : 3 Sep 2019 17:51
 Operator : HP/JU
 Sample : K4622-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG082219.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.115	3	4	16	rVB	117149	128344	7.82%	1.339%
2	5.553	413	419	433	rBV	319380	521300	31.76%	5.440%
3	7.163	686	693	709	rBV	320095	609656	37.15%	6.363%
4	7.527	748	755	771	rBV	353801	616608	37.57%	6.435%
5	7.997	825	835	845	rBV	103293	168942	10.29%	1.763%
6	8.321	881	890	900	rBV	285920	495156	30.17%	5.168%
7	9.167	1025	1034	1050	rBV	183455	359929	21.93%	3.756%
8	10.818	1308	1315	1334	rBV	119338	239113	14.57%	2.495%
9	13.274	1726	1733	1746	rBV	616374	1018080	62.03%	10.625%
10	14.108	1869	1875	1888	rBV	79394	137464	8.38%	1.435%
11	14.654	1962	1968	1982	rBV	237323	400831	24.42%	4.183%
12	16.147	2215	2222	2250	rBV	544038	959084	58.44%	10.009%
13	17.410	2430	2437	2450	rBV2	329901	531564	32.39%	5.548%
14	18.303	2585	2589	2593	rBV5	10682	18295	1.11%	0.191%
15	20.019	2875	2881	2894	rBV	1244904	1641196	100.00%	17.128%
16	21.323	3098	3103	3112	rBV2	59622	121402	7.40%	1.267%
17	21.687	3154	3165	3181	rBV	388919	713393	43.47%	7.445%
18	22.480	3297	3300	3306	rBV4	11529	19987	1.22%	0.209%
19	23.180	3413	3419	3426	rBV2	19187	38070	2.32%	0.397%
20	23.990	3551	3557	3566	rVB4	20327	48901	2.98%	0.510%
21	24.931	3706	3717	3733	rBV2	237298	750693	45.74%	7.835%
22	26.088	3903	3914	3923	rBV7	12217	43846	2.67%	0.458%

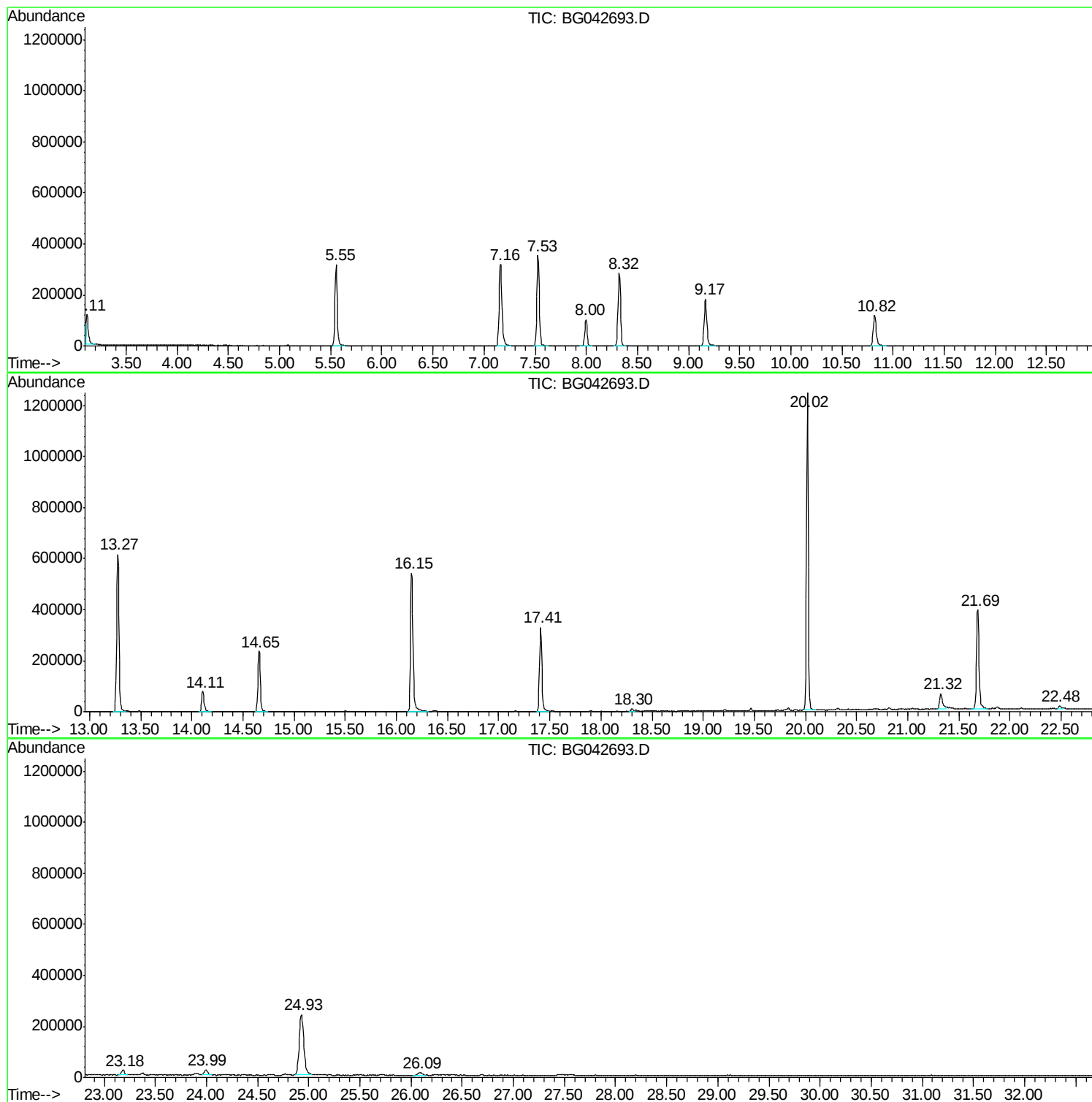
Sum of corrected areas: 9581854

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090319\
Data File : BG042693.D
Acq On : 3 Sep 2019 17:51
Operator : HP/JU
Sample : K4622-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090319\
Data File : BG042693.D
Acq On : 3 Sep 2019 17:51
Operator : HP/JU
Sample : K4622-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :

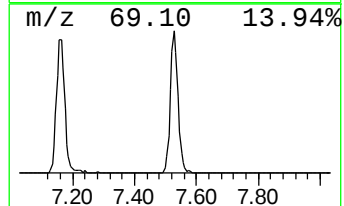
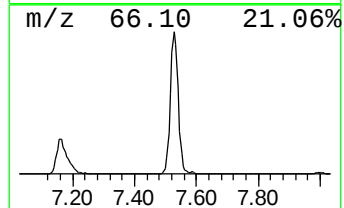
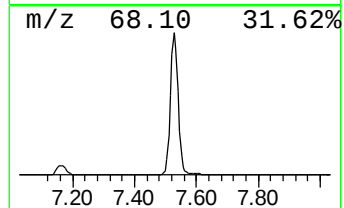
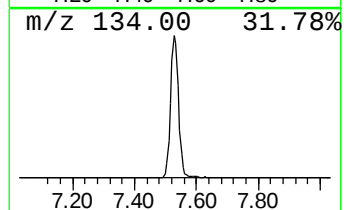
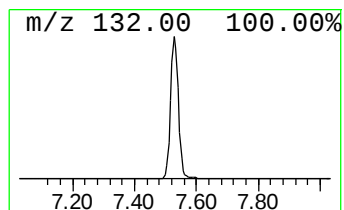
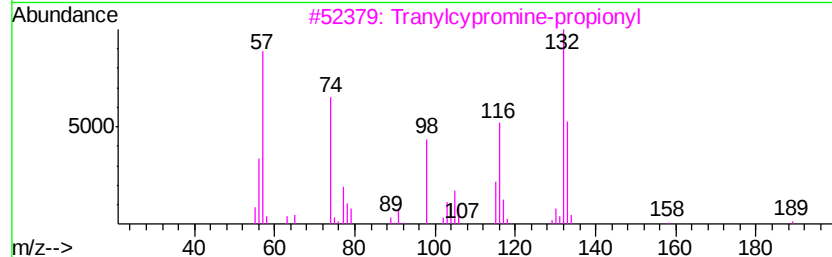
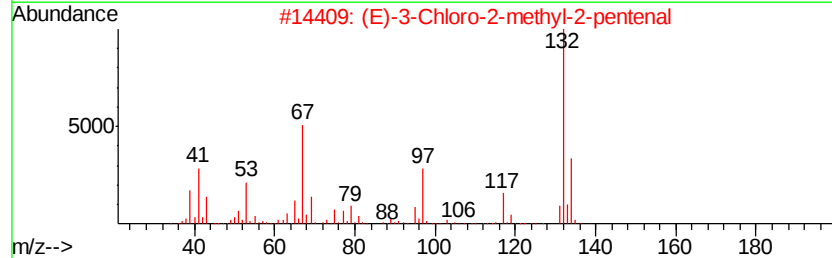
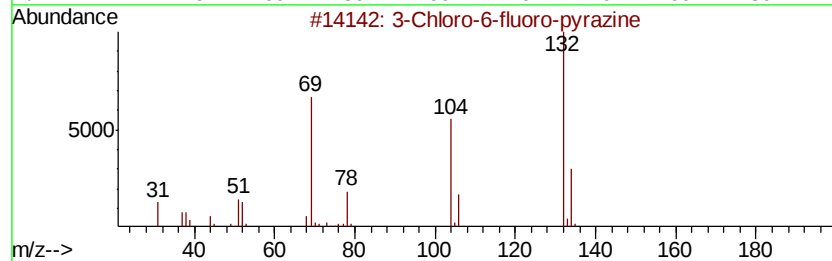
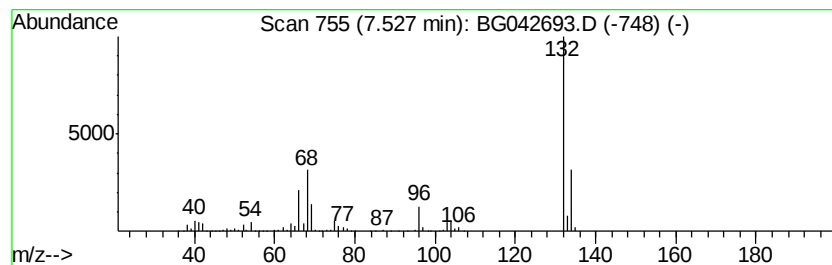
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	73.00 ng	616608	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	12
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090319\
Data File : BG042693.D
Acq On : 3 Sep 2019 17:51
Operator : HP/JU
Sample : K4622-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :

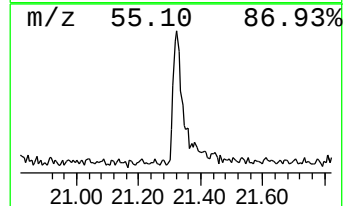
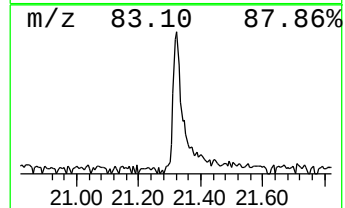
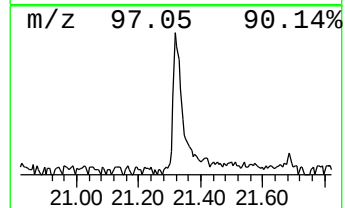
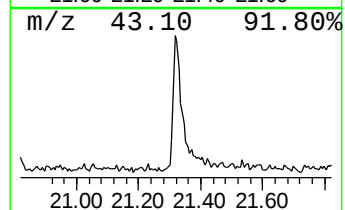
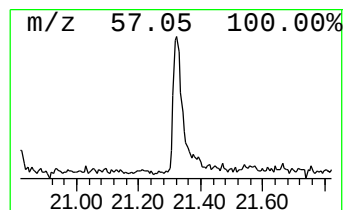
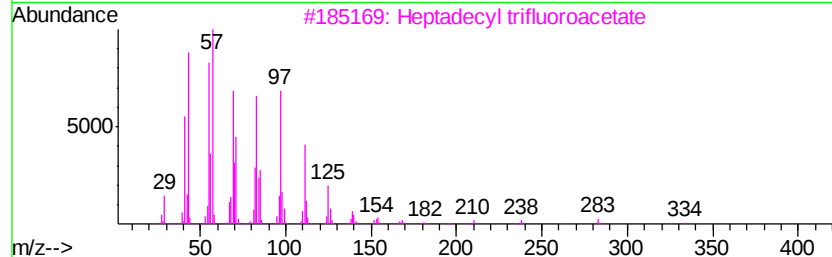
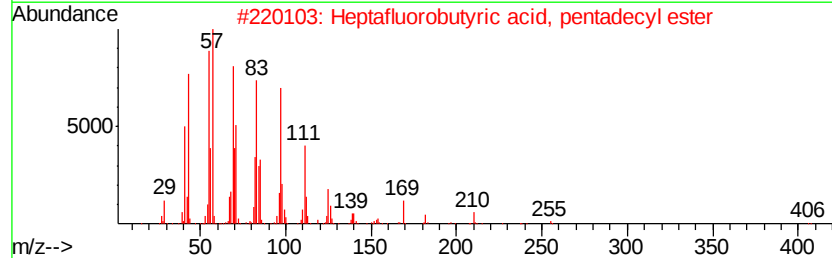
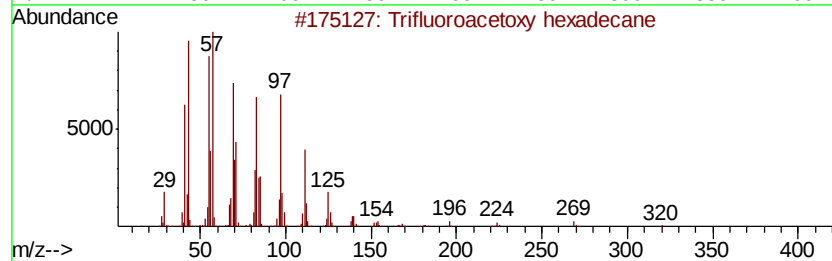
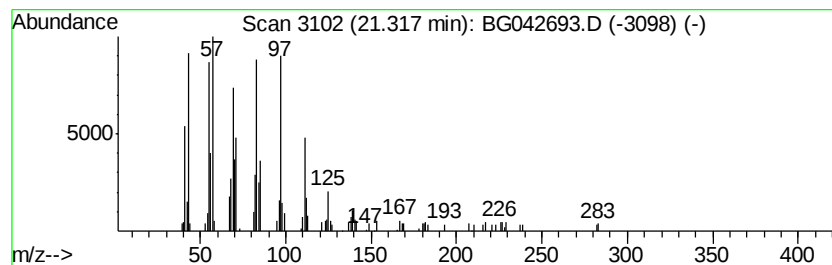
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Trifluoroacetoxy hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.32	3.40 ng	121402	Chrysene-d12	21.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
2		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
3		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
4		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	92
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG090319\
Data File : BG042693.D
Acq On : 3 Sep 2019 17:51
Operator : HP/JU
Sample : K4622-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.53	7.53	73.0	ng	616608	1	8.00	168942	20.0
Trifluoroacetoxy ...	21.32	3.4	ng	121402	5	21.68	713393	20.0