

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111822\
 Data File : BG055697.D
 Acq On : 18 Nov 2022 19:35
 Operator : CG/JU
 Sample : N5698-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BANK-RUN-SOIL-SAMPLE-1

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_G\Methods\8270-BG111622.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055697.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.296	337	342	349	rBV	89729	140544	4.86%	1.133%
2	5.778	415	424	433	rBV	563817	894723	30.95%	7.216%
3	7.394	690	699	709	rBV	570984	970115	33.55%	7.824%
4	8.246	838	844	851	rBV	98264	164843	5.70%	1.329%
5	9.421	1035	1044	1053	rBV	327998	591033	20.44%	4.766%
6	11.078	1318	1326	1337	rBV	139305	252050	8.72%	2.033%
7	13.499	1730	1738	1745	rBV	1014628	1597870	55.27%	12.886%
8	14.868	1965	1971	1980	rVB	271391	435145	15.05%	3.509%
9	16.213	2194	2200	2207	rBV	262168	370438	12.81%	2.987%
10	16.354	2217	2224	2239	rBV	1051410	1691761	58.51%	13.643%
11	16.542	2250	2256	2262	rBV	44430	67312	2.33%	0.543%
12	17.611	2431	2438	2445	rBV	434579	643243	22.25%	5.187%
13	20.202	2873	2879	2885	rBV	2159998	2891218	100.00%	23.317%
14	21.513	3097	3102	3115	rBV2	99911	247123	8.55%	1.993%
15	21.912	3163	3170	3176	rBV2	370718	649589	22.47%	5.239%
16	22.083	3196	3199	3203	rVB2	23022	32352	1.12%	0.261%
17	22.717	3304	3307	3313	rBV2	21906	35293	1.22%	0.285%
18	25.320	3741	3750	3762	rVB2	245826	725225	25.08%	5.849%

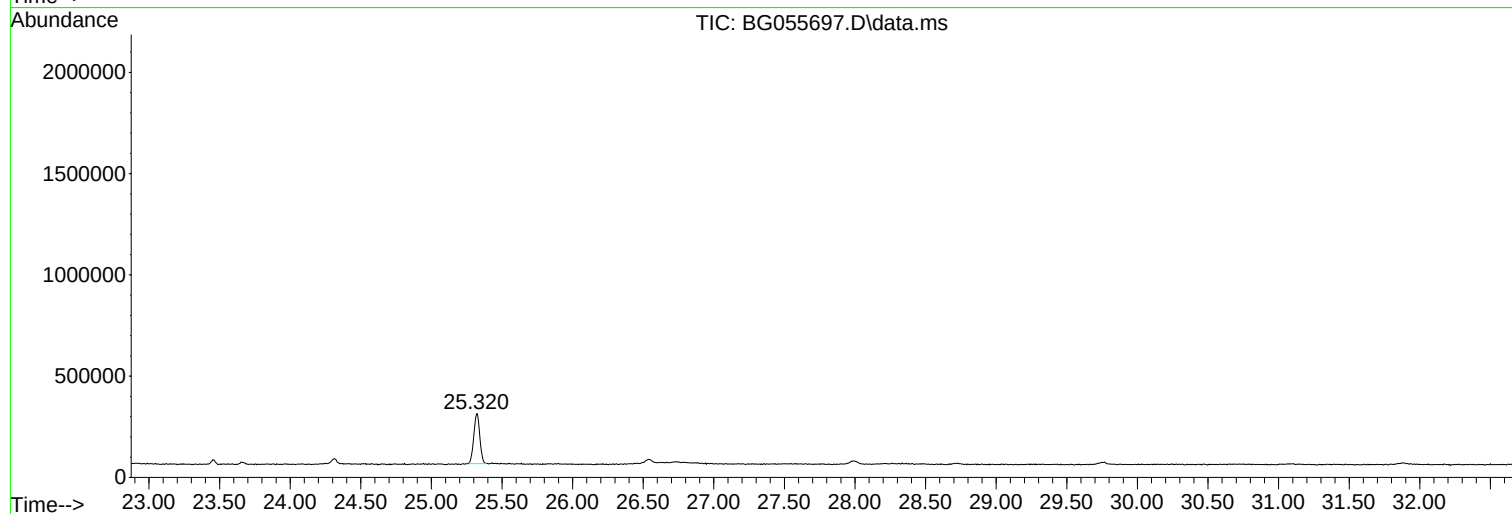
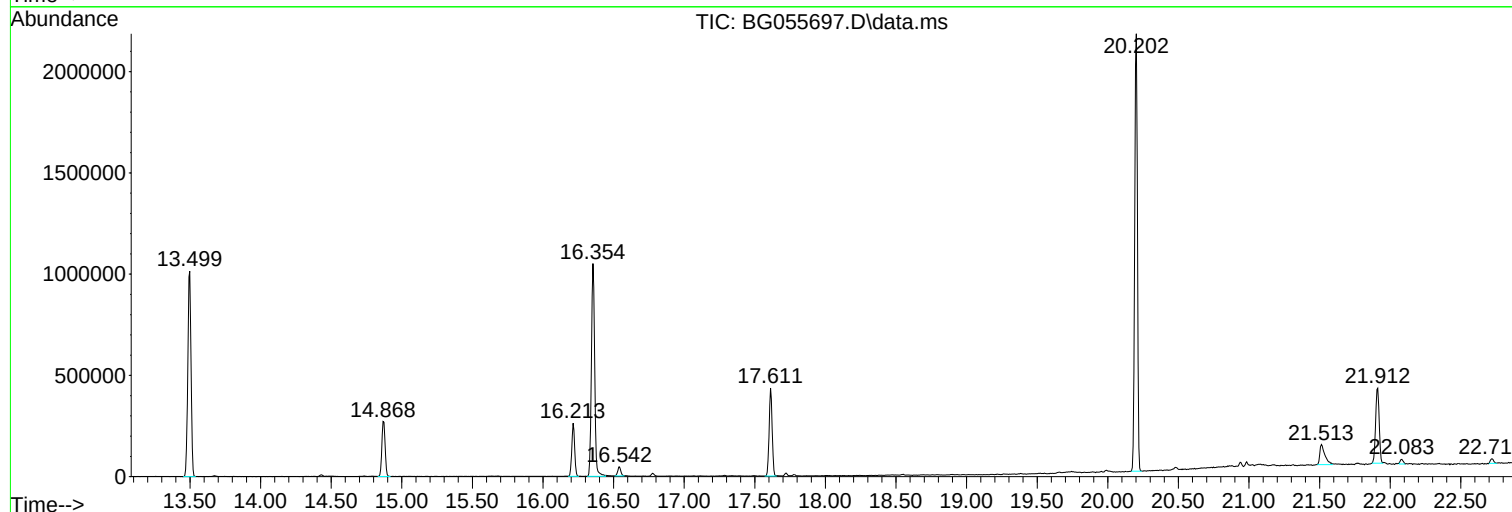
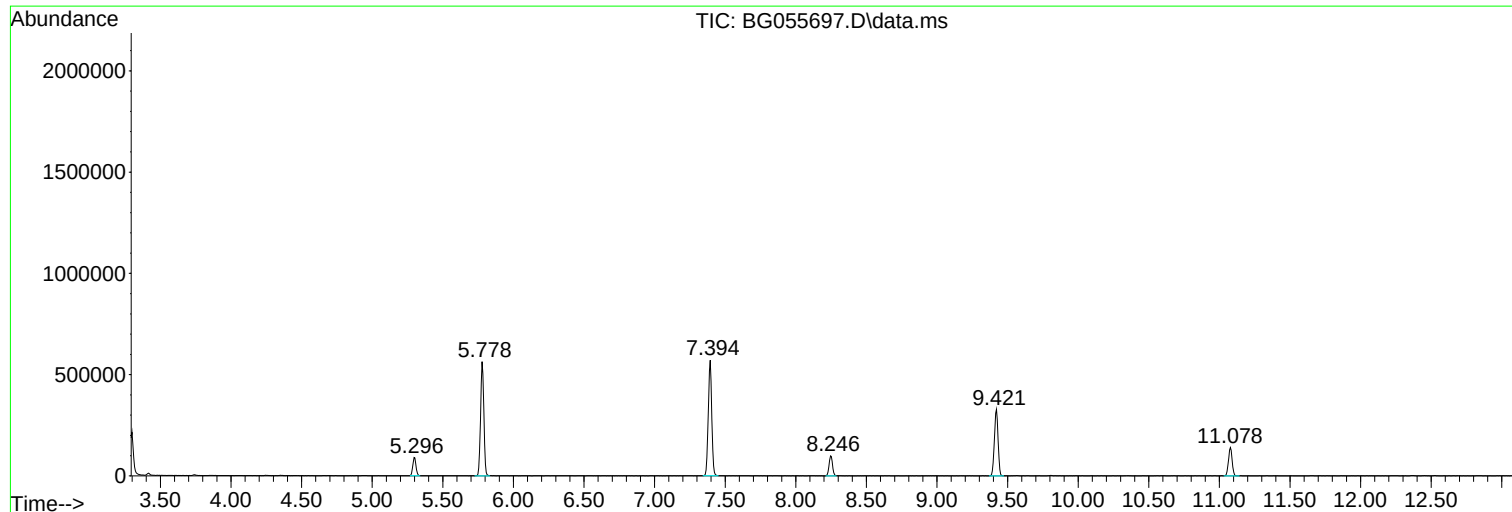
Sum of corrected areas: 12399877

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.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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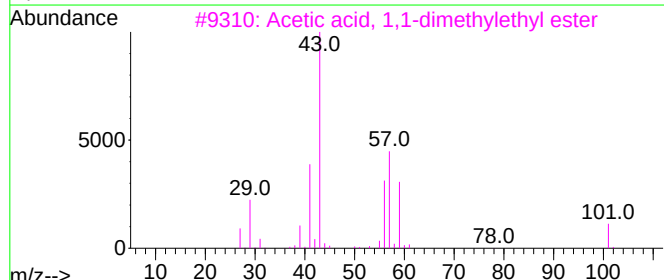
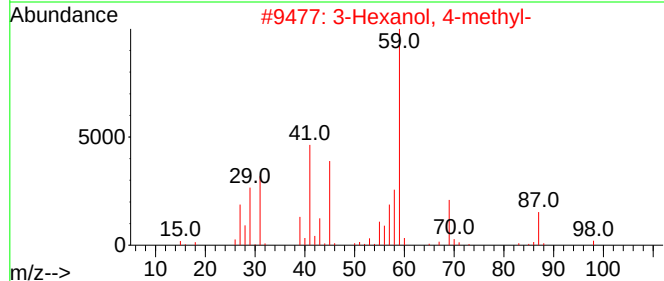
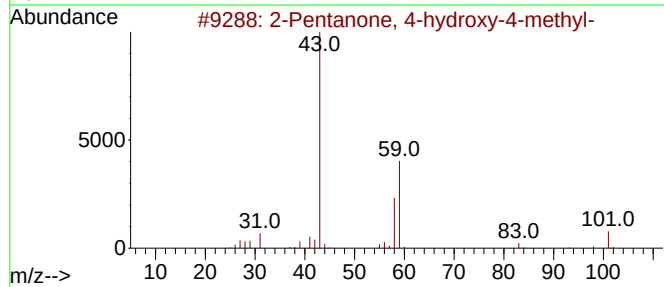
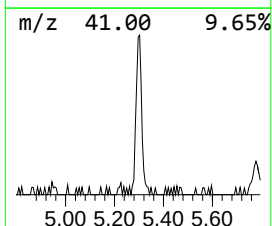
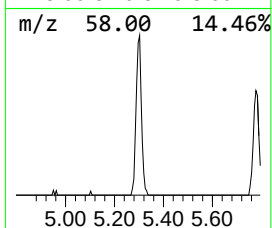
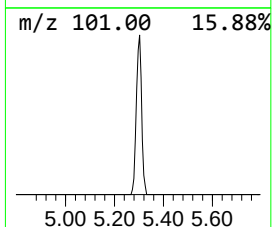
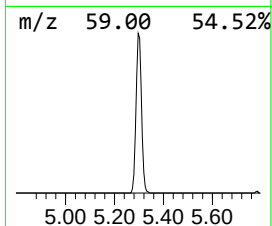
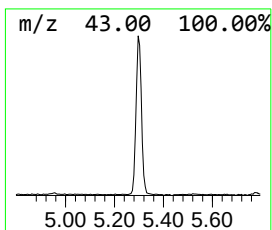
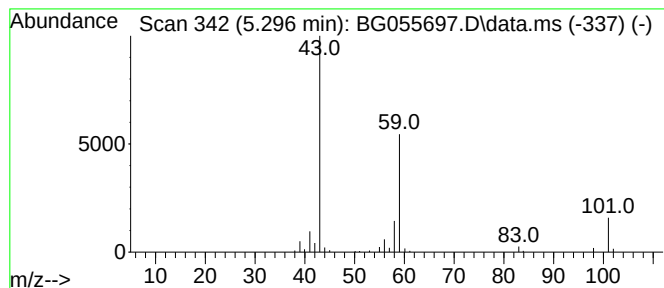
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.296	17.05 ng	140544	1,4-Dichlorobenzene-d4	8.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	32		
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		



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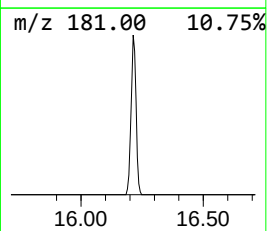
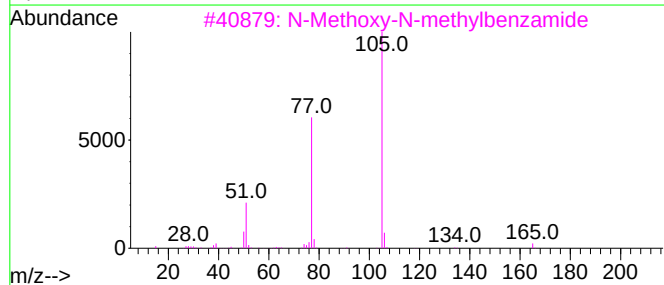
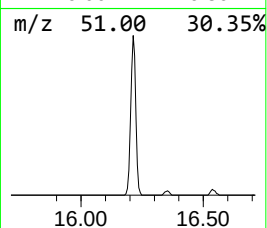
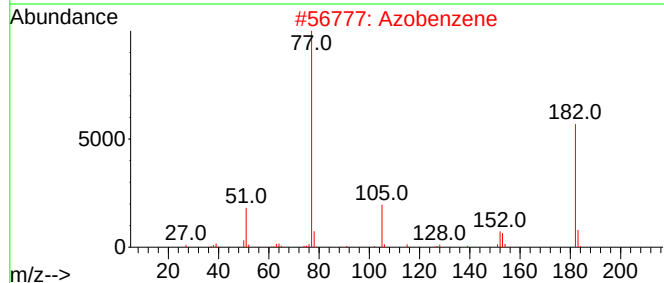
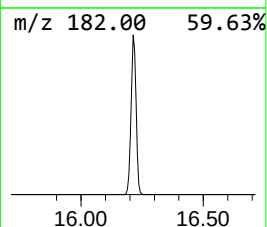
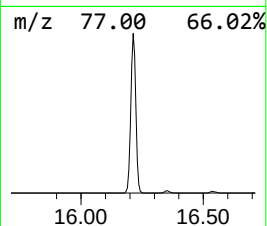
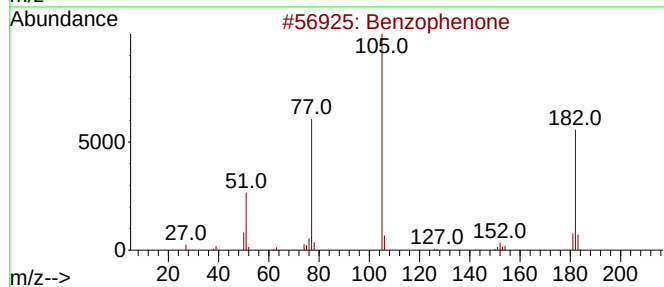
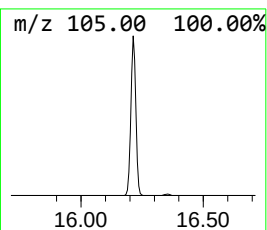
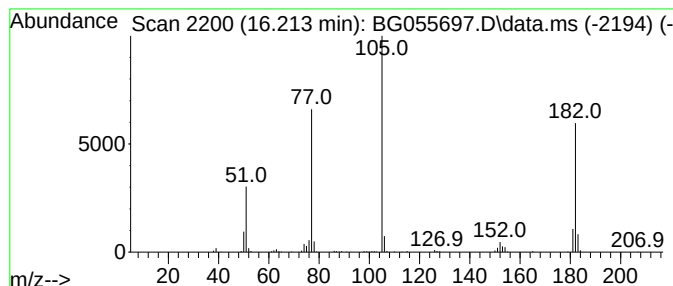
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.213	17.03 ng	370438	Acenaphthene-d10	14.868

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Azobenzene	182	C12H10N2	000103-33-3	52
3			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	47
4			Ethanedione, diphenyl-	210	C14H10O2	000134-81-6	47
5			Pentafluoroethyl phenyl ketone	224	C9H5F5O	000394-52-5	47



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Instrument :
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ClientSampleId :
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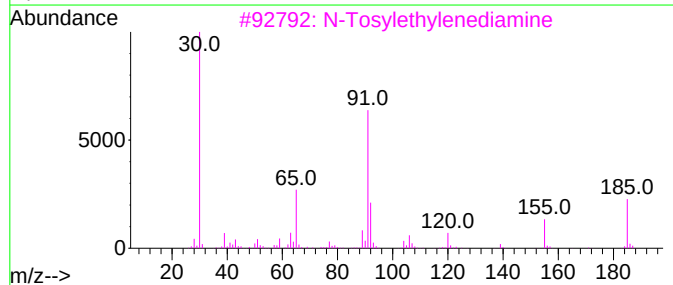
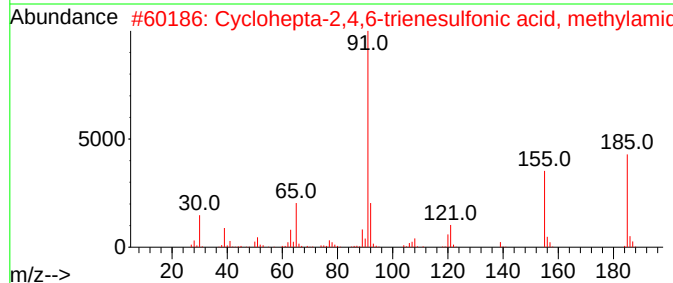
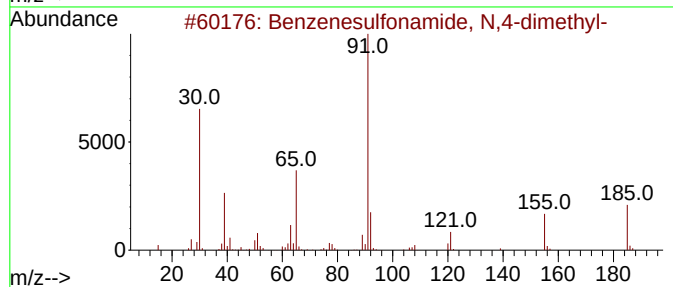
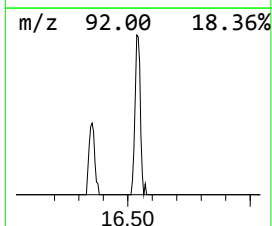
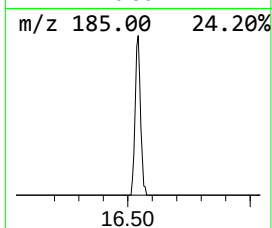
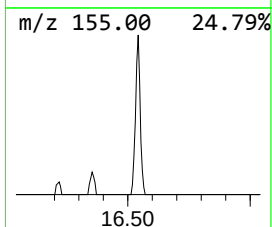
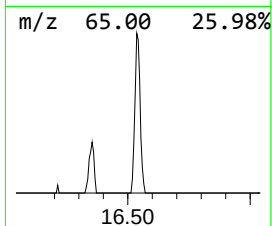
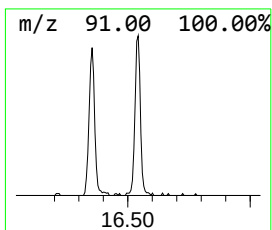
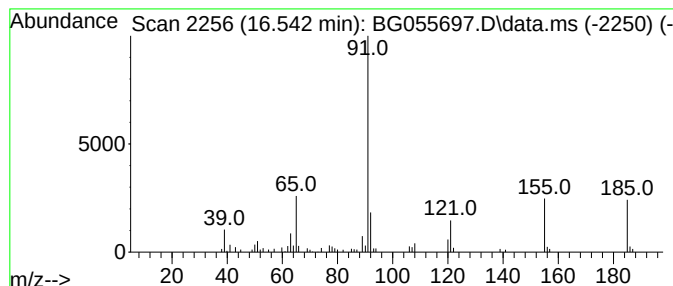
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Peak Number 3 Benzenesulfonamide, N,4-dim... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.542	2.09 ng	67312	Phenanthrene-d10	17.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	95
2			Cyclohepta-2,4,6-trienesulfonic ...	185	C8H11NO2S	1000187-28-7	76
3			N-Tosylethylenediamine	214	C9H14N2O2S	014316-16-6	53
4			Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	50
5			Pyridine-2-carbonitrile, 3,5,6-t...	375	C13H8Cl3N3O2S	255713-77-0	43



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ClientSampleId :
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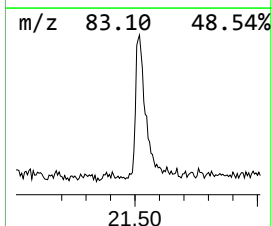
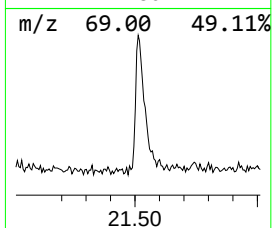
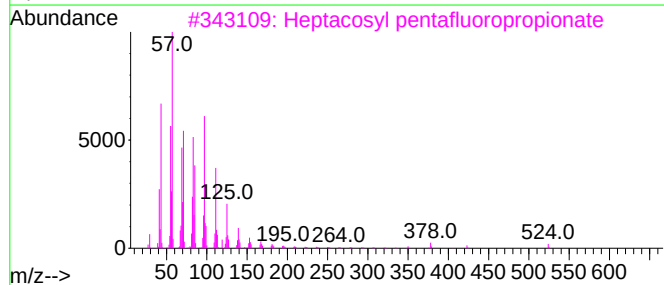
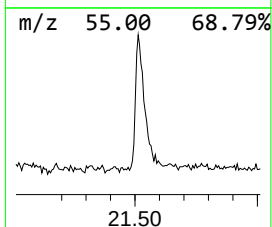
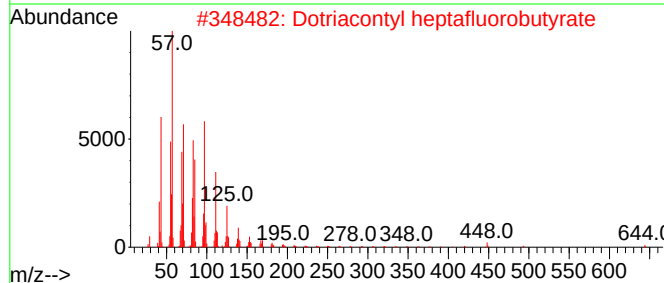
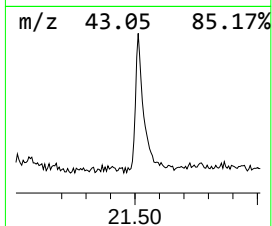
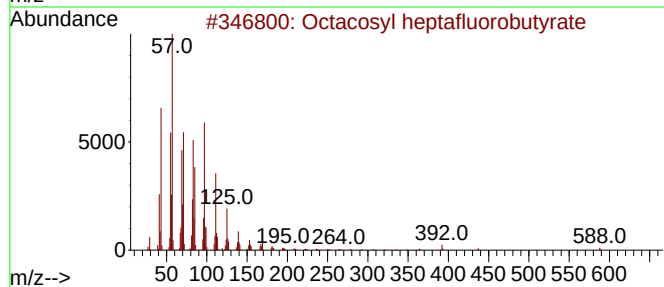
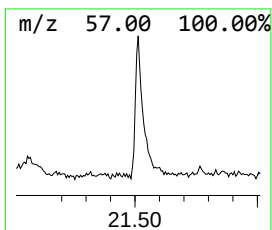
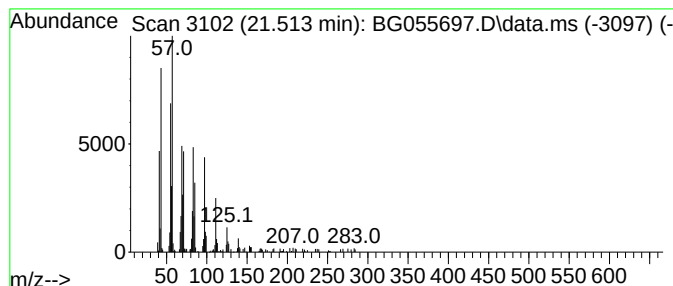
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Octacosyl heptafluorobutyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.513	7.61 ng	247123	Chrysene-d12	21.912

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosyl heptafluorobutyrate	606	C32H57F7O2	1010351-83-6	91
2			Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	91
3			Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	91
4			Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	91
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.296	17.1	ng	140544	1	8.246	164843	20.0
Benzophenone	16.213	17.0	ng	370438	3	14.868	435145	20.0
Benzenesulfonam...	16.542	2.1	ng	67312	4	17.611	643243	20.0
Octacosyl hepta...	21.513	7.6	ng	247123	5	21.912	649589	20.0