

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062618\
 Data File : BF106861.D
 Acq On : 27 Jun 2018 9:03
 Operator : JU/SJ
 Sample : J3693-02
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-26

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-BF061918.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	5.189	481	485	494	rBB	34311	39400	2.95%	0.472%
2	5.595	551	554	567	rBV	511724	480987	36.05%	5.764%
3	6.595	721	724	741	rBB	327421	291474	21.85%	3.493%
4	6.736	743	748	751	rBV	912725	843440	63.22%	10.108%
5	6.954	782	785	789	rBV	272156	244374	18.32%	2.929%
6	7.113	808	812	815	rBV	858229	762063	57.12%	9.133%
7	7.525	877	882	885	rBV	678579	654208	49.04%	7.840%
8	8.242	1000	1004	1018	rBV	284499	276744	20.74%	3.317%
9	9.313	1182	1186	1195	rBV	1206114	1076603	80.70%	12.903%
10	9.701	1247	1252	1259	rVB	18634	17884	1.34%	0.214%
11	9.995	1299	1302	1306	rBV	285026	278928	20.91%	3.343%
12	10.030	1306	1308	1311	rVB	23177	16676	1.25%	0.200%
13	10.654	1410	1414	1417	rBV	267258	244797	18.35%	2.934%
14	10.795	1434	1438	1441	rBV	757048	698013	52.32%	8.365%
15	10.848	1443	1447	1451	rBV	182068	171410	12.85%	2.054%
16	11.489	1552	1556	1560	rBV	358920	299436	22.45%	3.589%
17	12.871	1789	1791	1796	rVB3	19538	17156	1.29%	0.206%
18	13.077	1821	1826	1829	rBV	1414057	1334068	100.00%	15.988%
19	13.954	1971	1975	1983	rVB	32868	39072	2.93%	0.468%
20	14.142	2003	2007	2014	rBV	374213	344292	25.81%	4.126%
21	15.677	2261	2268	2275	rVB	154002	213007	15.97%	2.553%

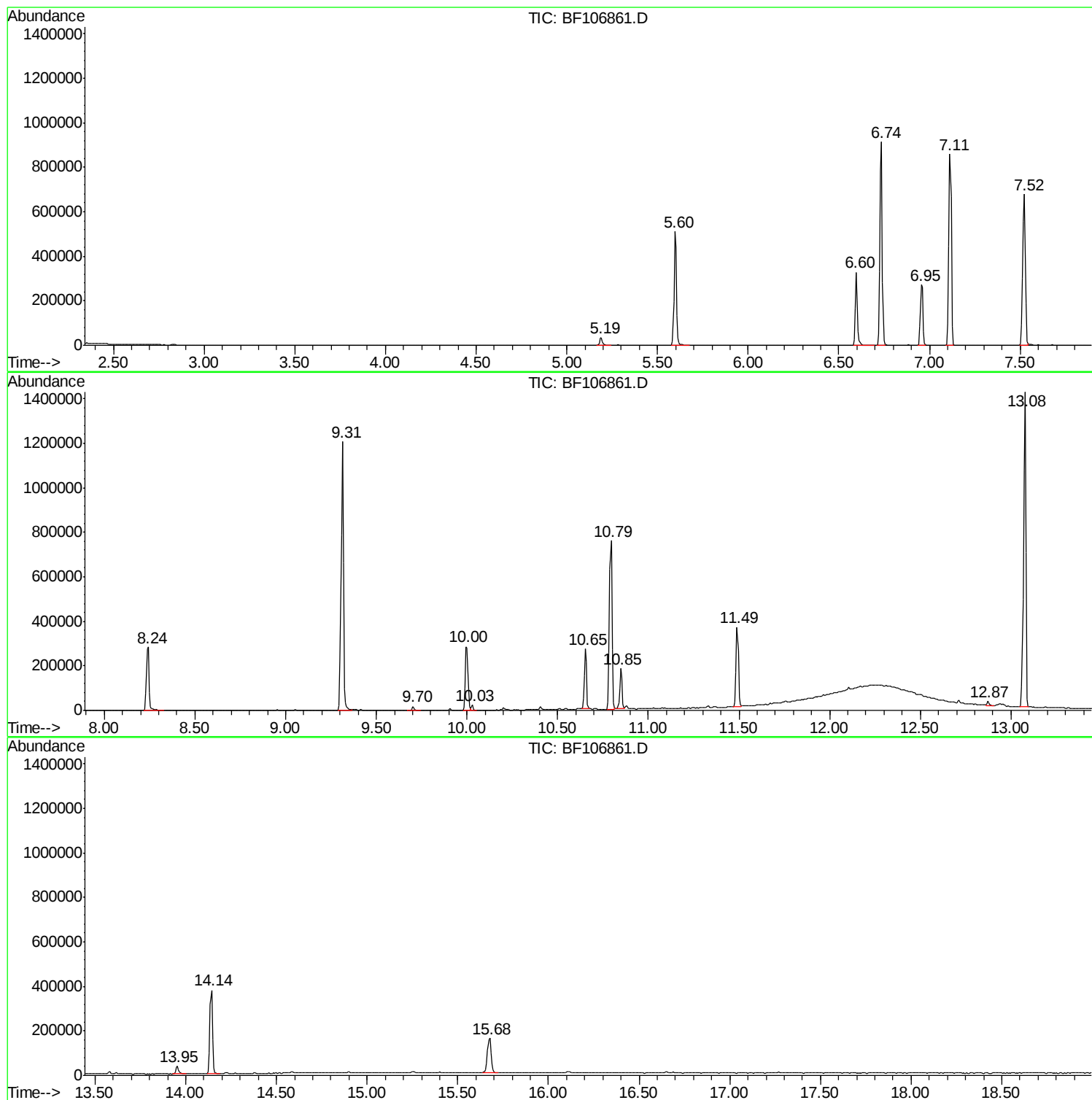
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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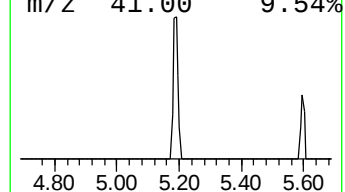
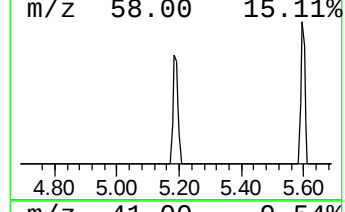
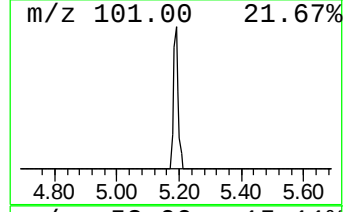
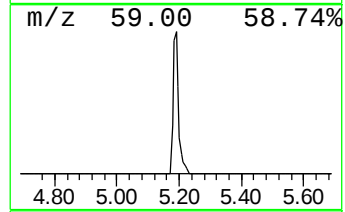
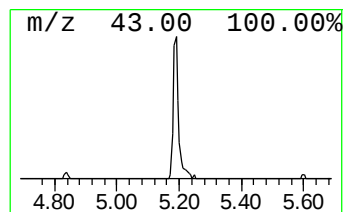
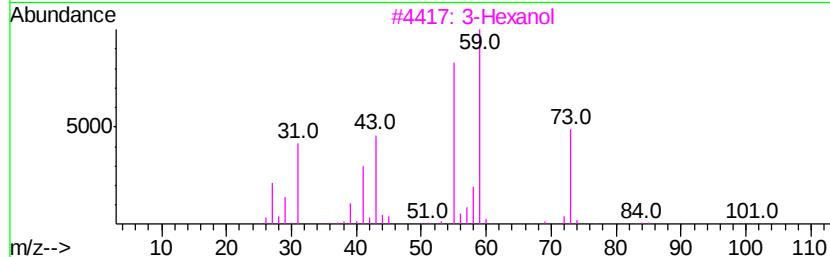
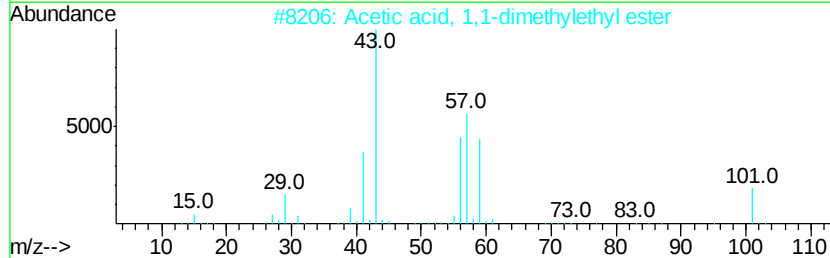
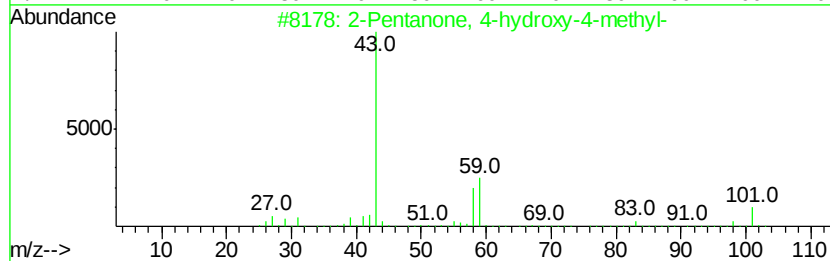
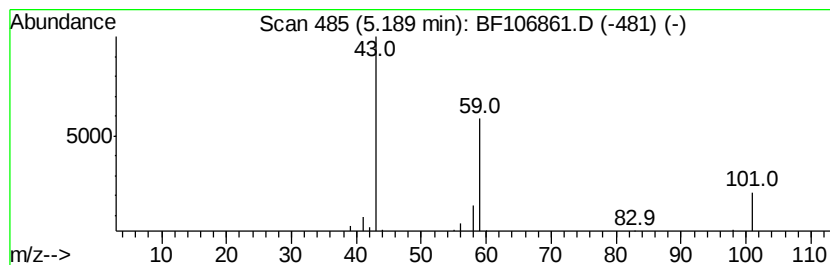
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.22 ng	39400	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	40
3			3-Hexanol	102	C6H14O	000623-37-0	33
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
5			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	9



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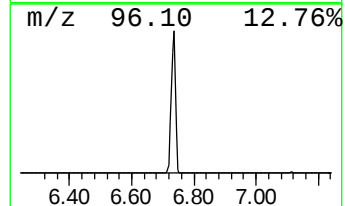
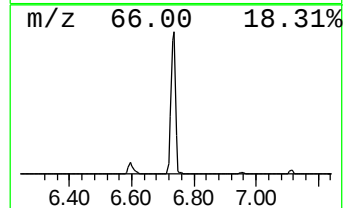
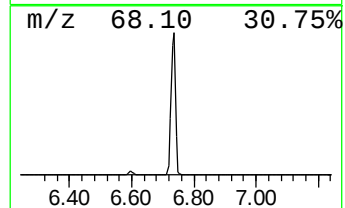
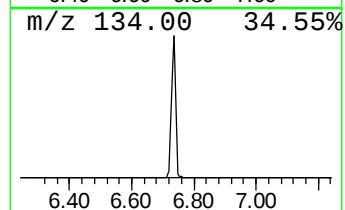
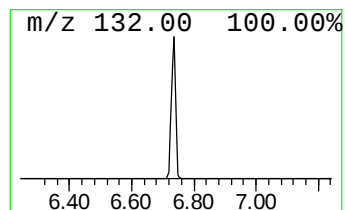
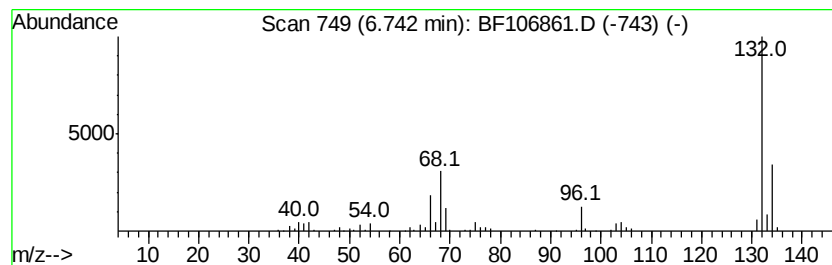
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	69.03 ng	843440	1,4-Dichlorobenzene-d4	6.96

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	9



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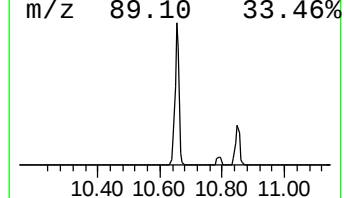
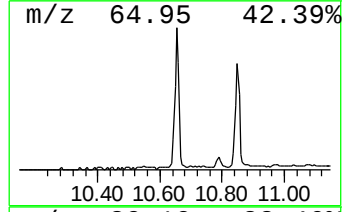
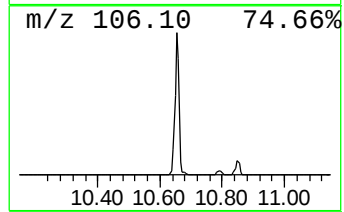
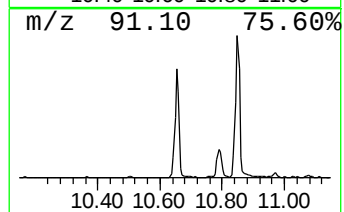
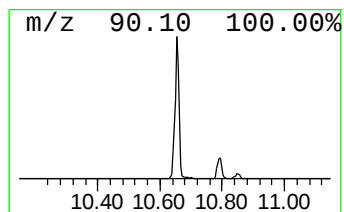
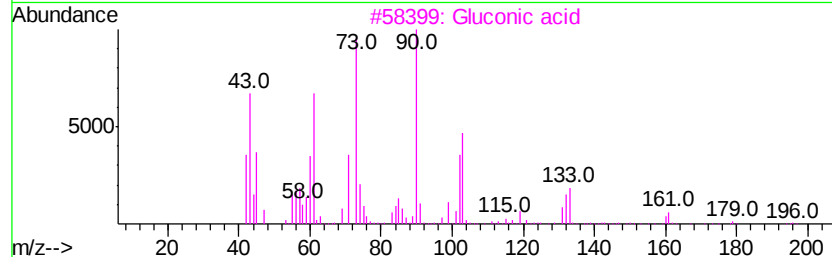
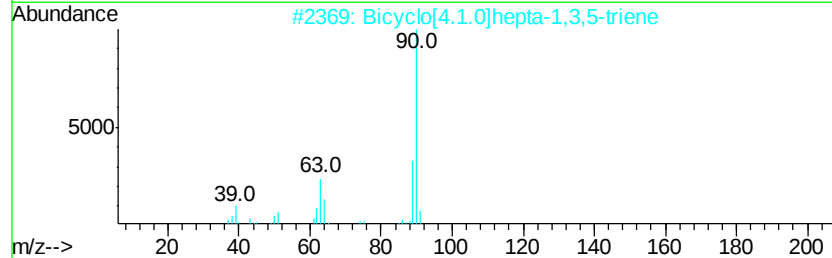
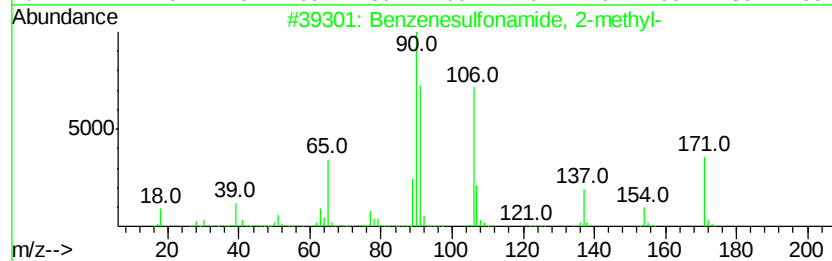
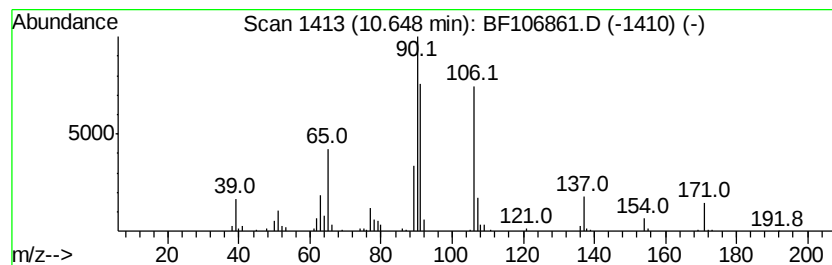
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Peak Number 4 Benzenesulfonamide, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	17.55 ng	244797	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, 2-methyl-	171	C7H9NO2S	000088-19-7	96
2		Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	46
3		Gluconic acid	196	C6H12O7	000526-95-4	38
4		4-(2-Aminoethyl)benzenesulfonamide	200	C8H12N2O2S	035303-76-5	37
5		Benzenemethanesulfonamide	171	C7H9NO2S	004563-33-1	20



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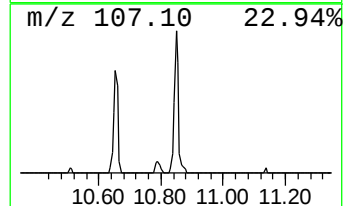
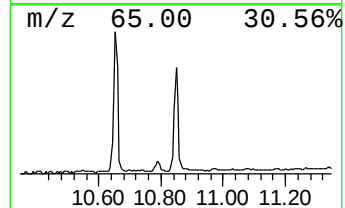
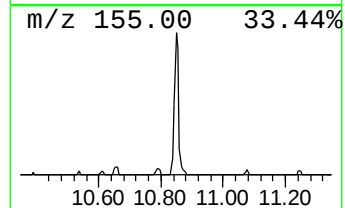
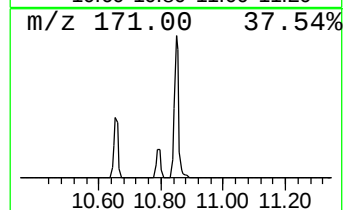
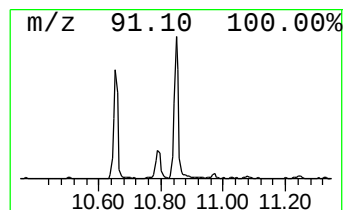
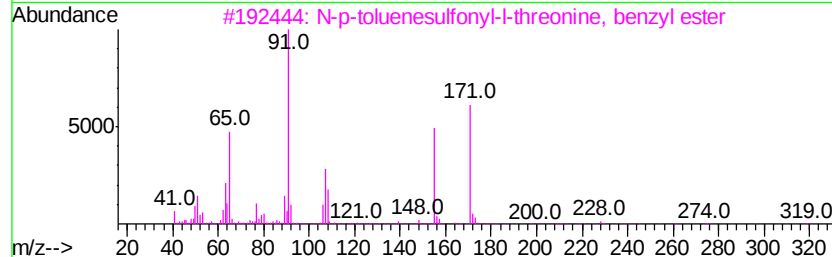
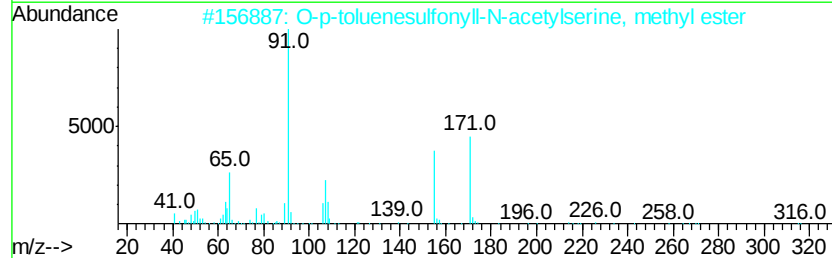
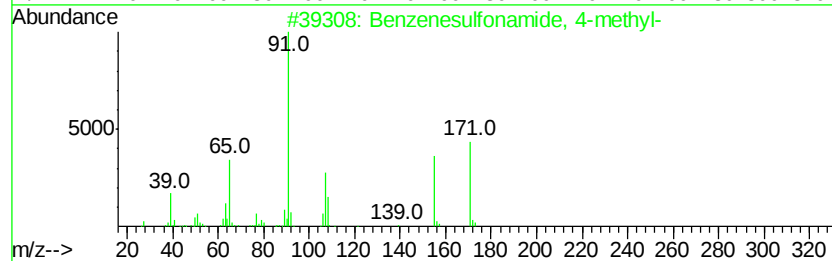
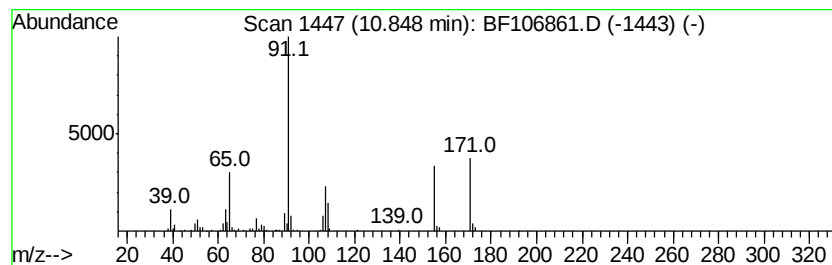
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Peak Number 5 Benzenesulfonamide, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	11.45 ng	171410	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, 4-methyl-	171	C7H9NO2S	000070-55-3	98
2		O-p-toluenesulfonyl-N-acetylser...	315	C13H17NO6S	1000126-44-8	90
3		N-p-toluenesulfonyl-L-threonine,...	363	C18H21NO5S	1000130-34-4	80
4		Benzenemethanesulfonamide	171	C7H9NO2S	004563-33-1	43
5		Pyridine-2-carbonitrile, 3,5,6-t...	375	C13H8Cl3N3O2S	255713-77-0	43



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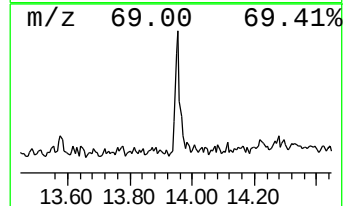
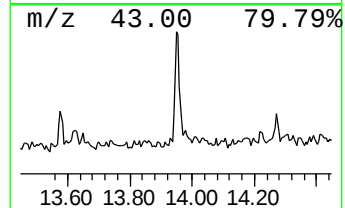
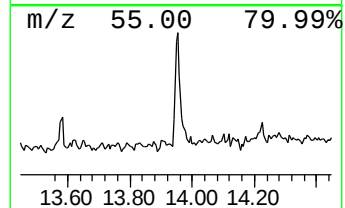
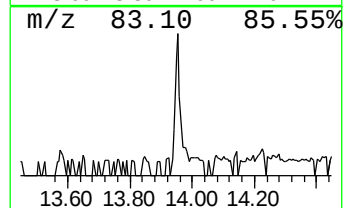
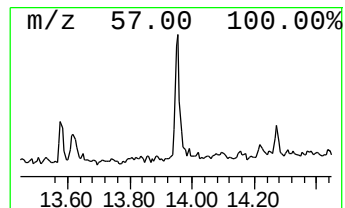
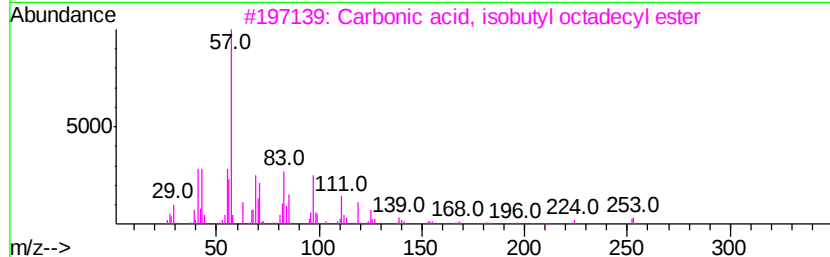
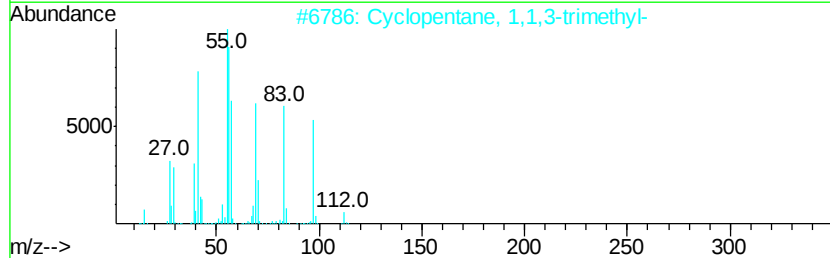
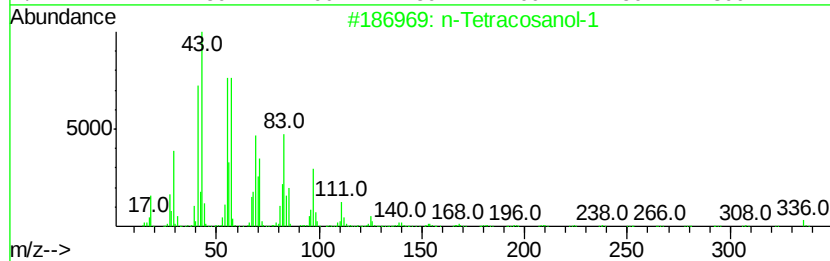
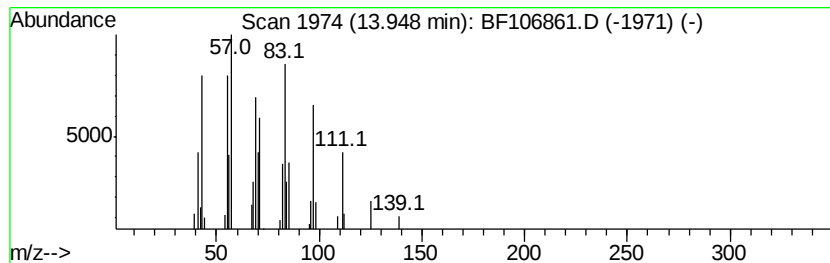
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Peak Number 6 n-Tetracosanol-1 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	2.27 ng	39072	Chrysene-d12	14.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		n-Tetracosanol-1	354 C24H50O	000506-51-4 83
2		Cyclopentane, 1,1,3-trimethyl-	112 C8H16	004516-69-2 49
3		Carbonic acid, isobutyl octadecyl...	370 C23H46O3	1000314-61-5 49
4		Isobutyl tetradecyl carbonate	314 C19H38O3	959275-58-2 47
5		Cycloheptane, methyl-	112 C8H16	004126-78-7 41



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.2	ng	39400	1	6.96	244374	20.0
unknown6.74	6.74	69.0	ng	843440	1	6.96	244374	20.0
Benzenesulfonamid...	10.65	17.6	ng	244797	3	10.00	278928	20.0
Benzenesulfonamid...	10.85	11.4	ng	171410	4	11.49	299436	20.0
n-Tetracosanol-1	13.95	2.3	ng	39072	5	14.14	344292	20.0