

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF120117\
 Data File : BF101060.D
 Acq On : 1 Dec 2017 13:57
 Operator : SJ/JU
 Sample : PB104640BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB104640BL

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:\HPCHEM1\BNA_F\METHODS\8270-BF113017.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	5.075	505	508	518	rBB	47977	63348	3.45%	0.518%
2	5.469	570	575	578	rBV	1219037	1207339	65.69%	9.876%
3	6.481	741	747	750	rBV	1217901	1232326	67.05%	10.080%
4	6.616	765	770	773	rBV	1440056	1289992	70.19%	10.552%
5	6.845	805	809	812	rBV	364411	289069	15.73%	2.365%
6	7.004	831	836	839	rBV	1077139	999034	54.36%	8.172%
7	7.410	900	905	908	rBV	826984	781720	42.53%	6.394%
8	8.128	1023	1027	1030	rBV	473696	386756	21.04%	3.164%
9	9.204	1205	1210	1213	rBV	1889238	1636483	89.04%	13.386%
10	9.592	1273	1276	1279	rBV	63858	47020	2.56%	0.385%
11	9.881	1321	1325	1329	rBV2	509384	444187	24.17%	3.633%
12	10.304	1394	1397	1400	rVB2	25748	20776	1.13%	0.170%
13	10.675	1456	1460	1463	rBV	694744	634360	34.52%	5.189%
14	10.775	1474	1477	1482	rBV	20211	19667	1.07%	0.161%
15	11.369	1574	1578	1582	rBV	562155	477980	26.01%	3.910%
16	12.963	1843	1849	1852	rBV	1807003	1837869	100.00%	15.034%
17	13.839	1995	1998	2002	rBV2	42254	47451	2.58%	0.388%
18	13.874	2002	2004	2008	rVB	34275	27178	1.48%	0.222%
19	14.016	2024	2028	2031	rBV	478652	437677	23.81%	3.580%
20	14.845	2165	2169	2175	rVB	30845	31460	1.71%	0.257%
21	15.480	2272	2277	2282	rBV	260201	313220	17.04%	2.562%

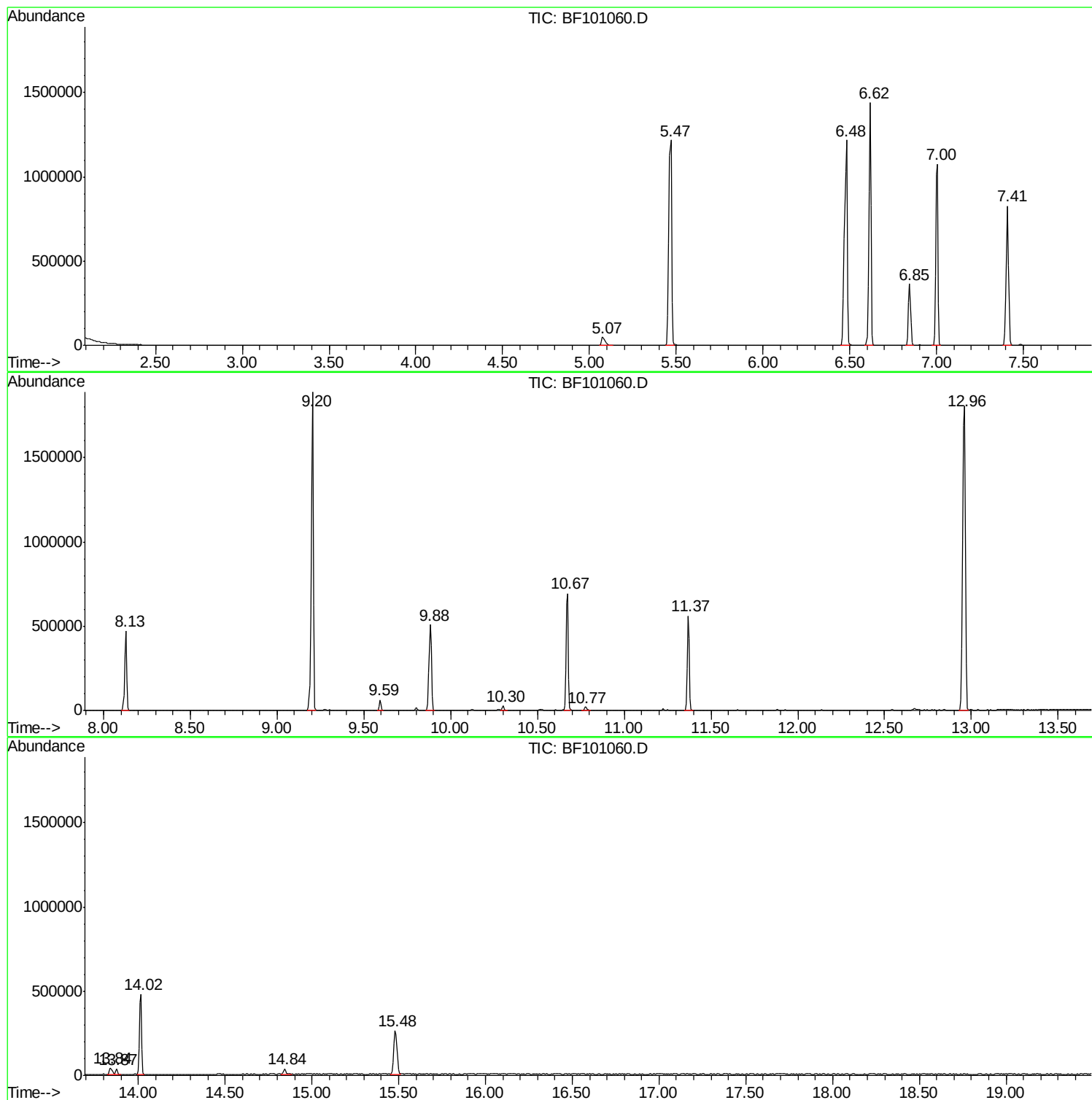
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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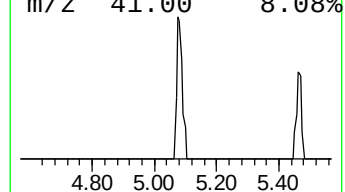
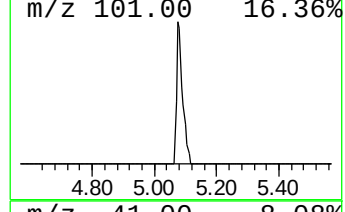
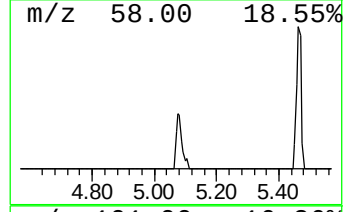
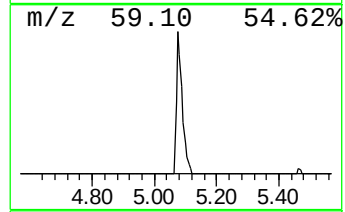
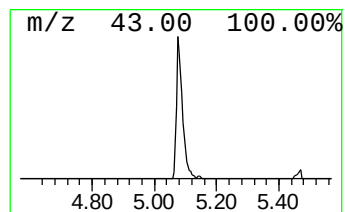
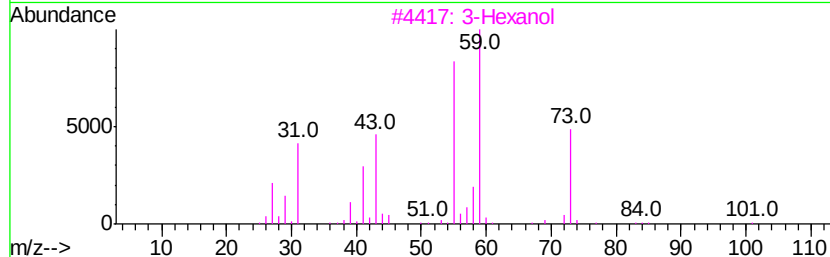
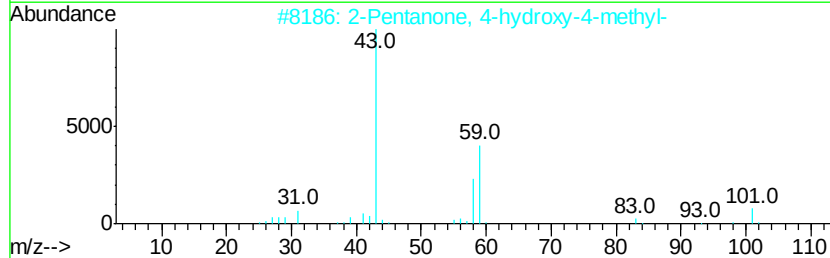
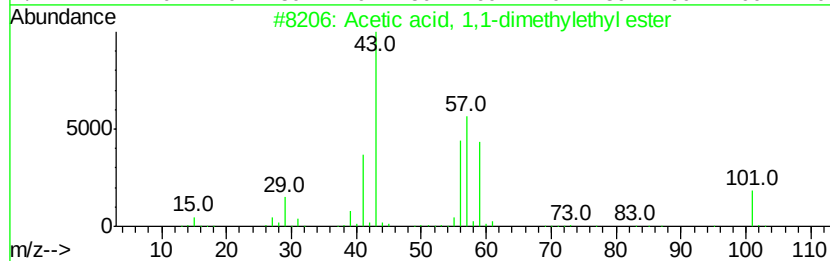
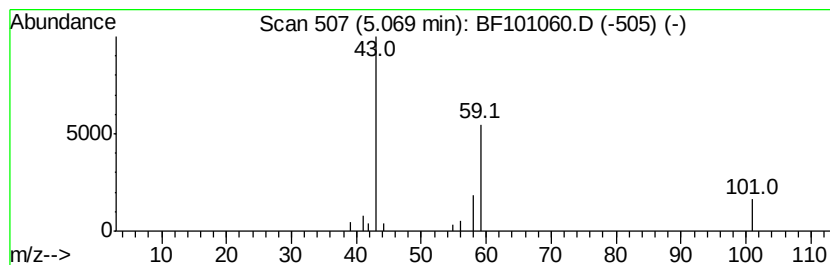
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown5.07 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	4.38 ng	63348	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
3			3-Hexanol	102	C6H14O	000623-37-0	36
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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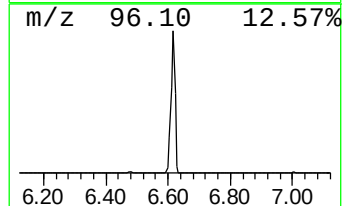
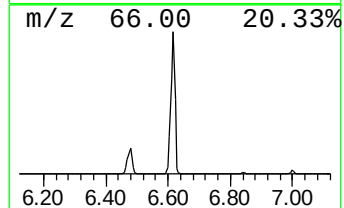
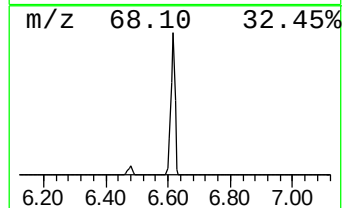
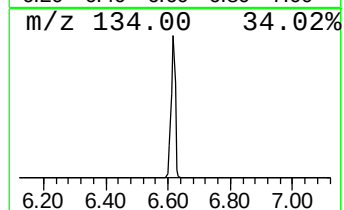
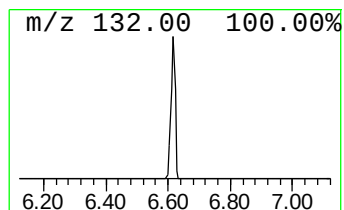
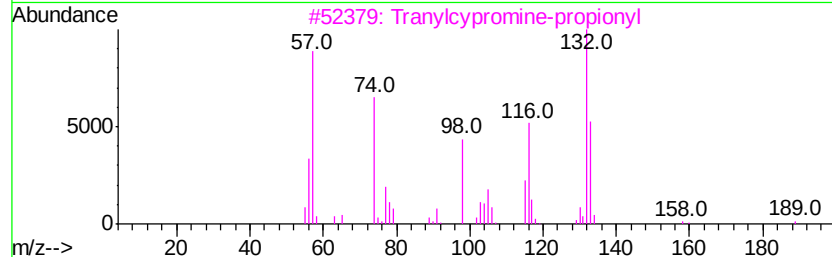
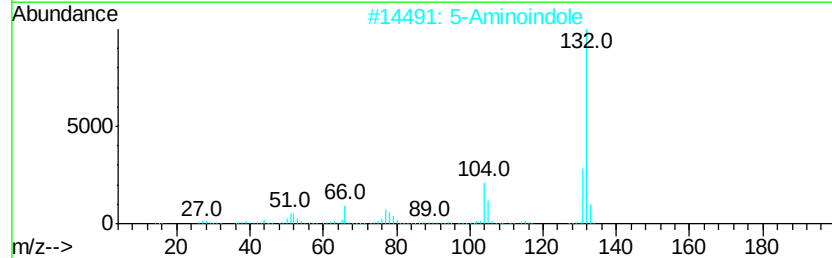
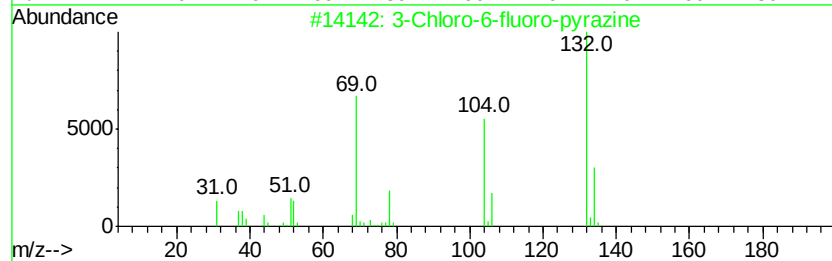
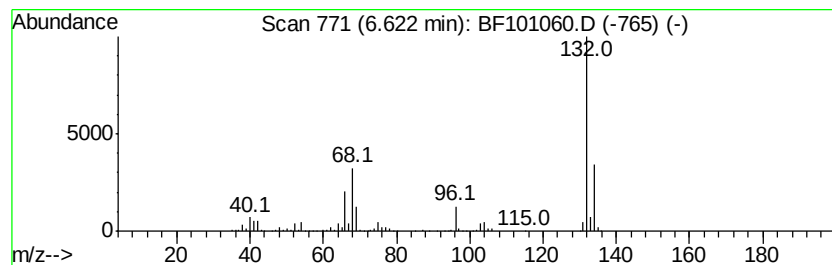
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	89.25 ng	1289990	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2	5		Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	10
4	4		Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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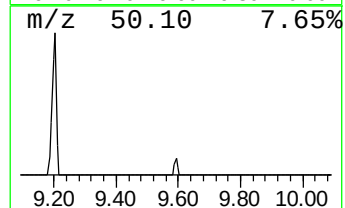
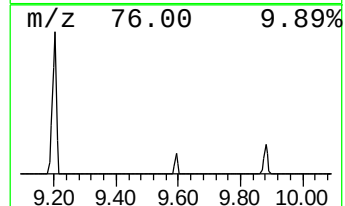
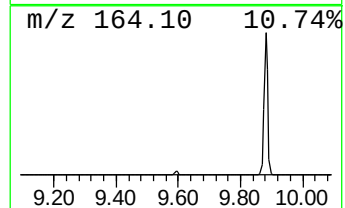
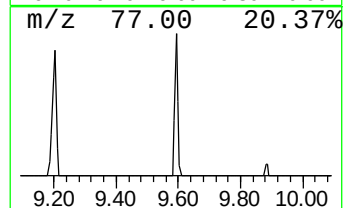
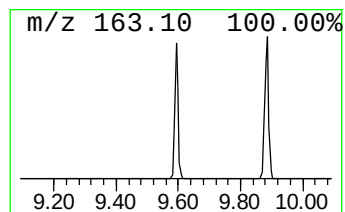
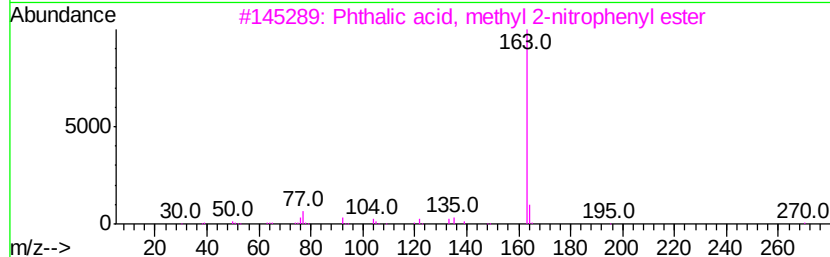
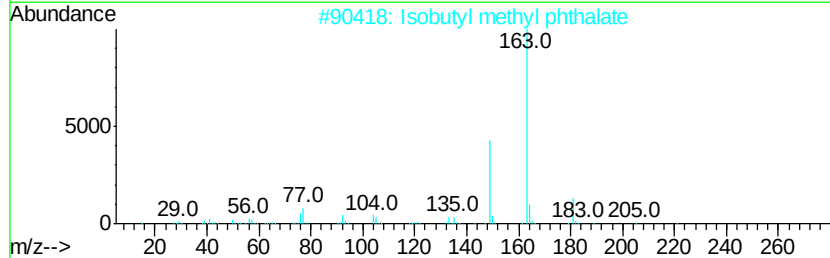
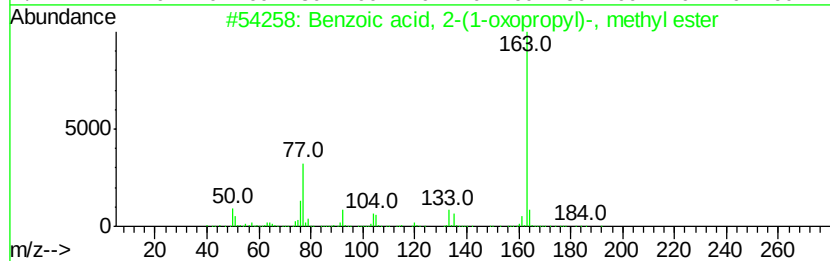
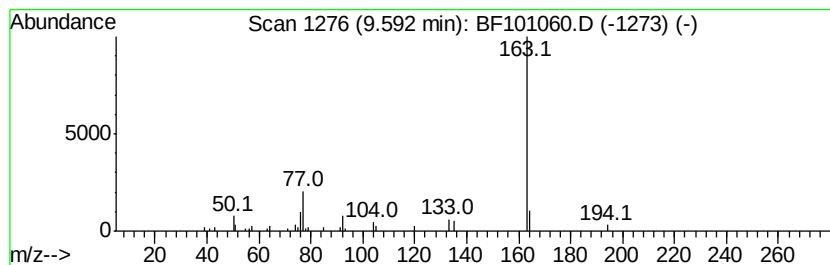
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Peak Number 4 Benzoic acid, 2-(1-oxopropyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.59	2.12 ng	47020	Acenaphthene-d10	9.88
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzoic acid, 2-(1-oxopropyl)-, ...	192 C11H12O3	032025-37-9 83
2		Isobutyl methyl phthalate	236 C13H16O4	1000373-89-3 78
3		Phthalic acid, methyl 2-nitrophenyl ester	301 C15H11NO6	1000315-68-3 74
4		Dimethyl phthalate	194 C10H10O4	000131-11-3 70
5		2,3-Dihydro-7-methyl-1,4-benzoxa...	163 C9H9NO2	039522-25-3 64



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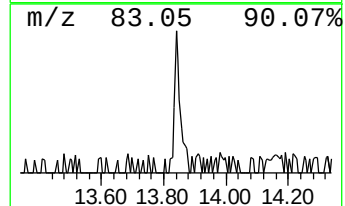
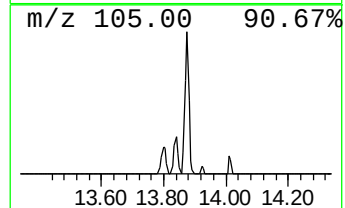
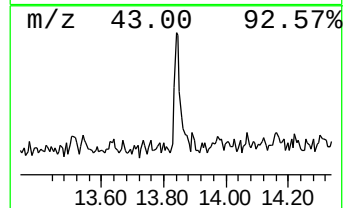
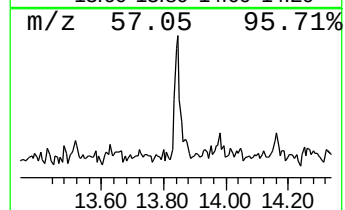
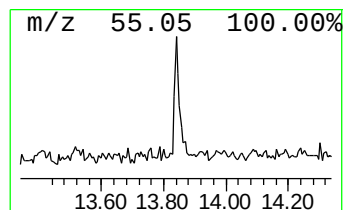
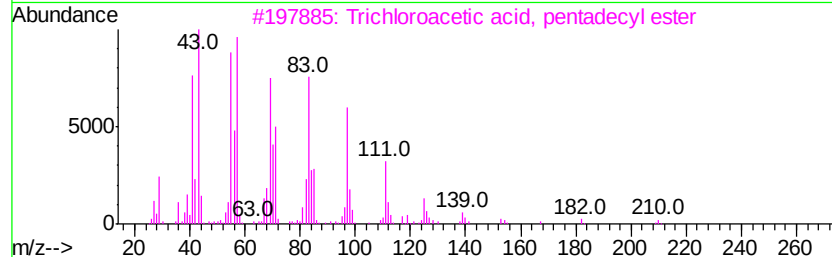
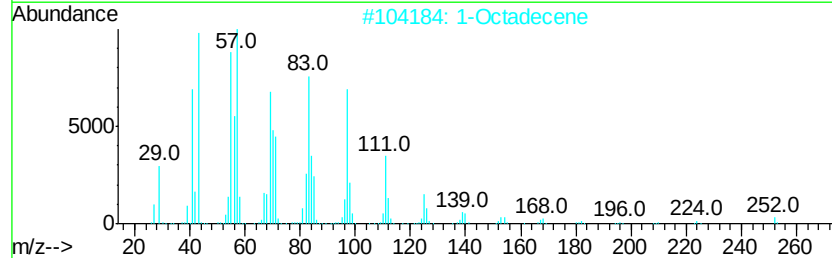
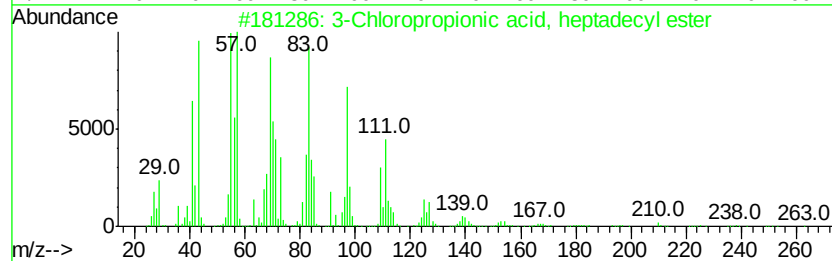
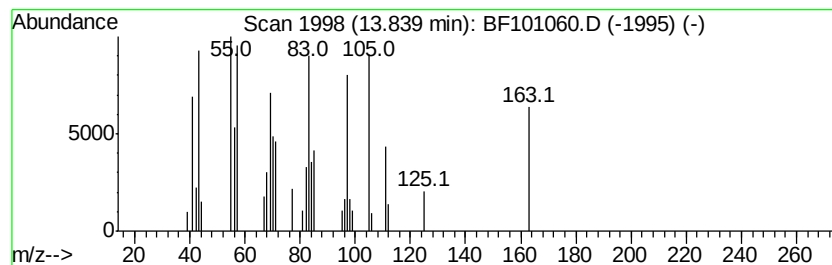
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Peak Number 5 3-Chloropropionic acid, hep... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	2.17 ng	47451	Chrysene-d12	14.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	87
2		1-Octadecene	252	C18H36	000112-88-9	87
3		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	87
4		11-Tricosene	322	C23H46	052078-56-5	87
5		1-Nonadecene	266	C19H38	018435-45-5	83



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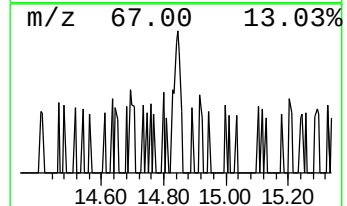
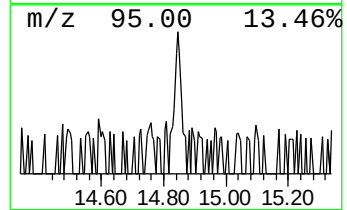
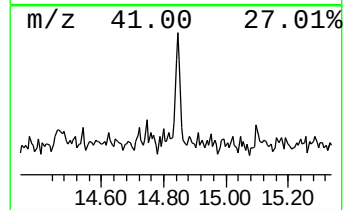
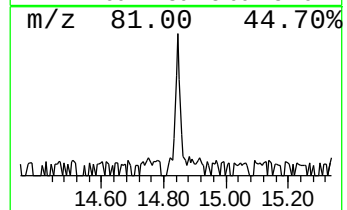
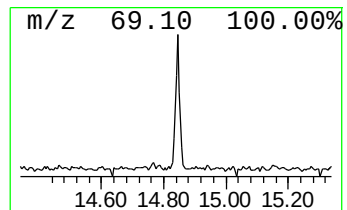
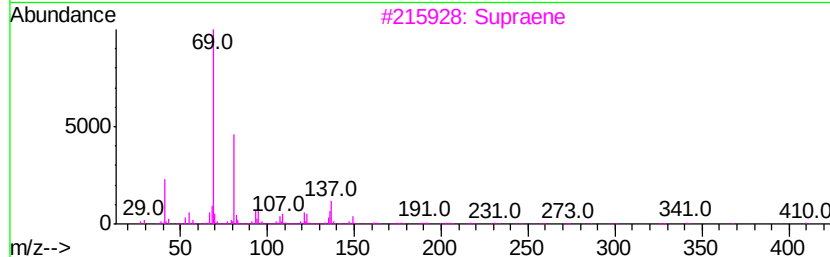
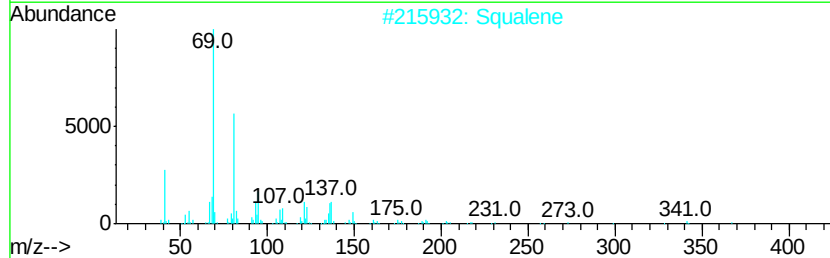
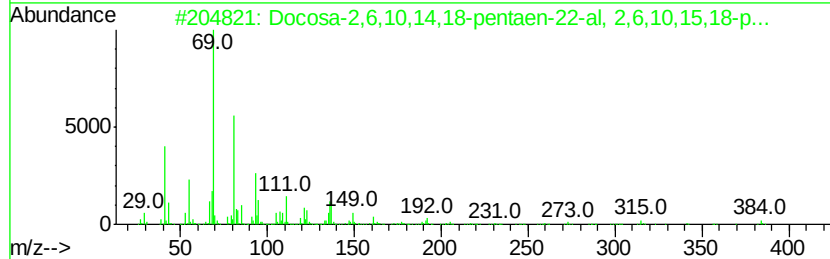
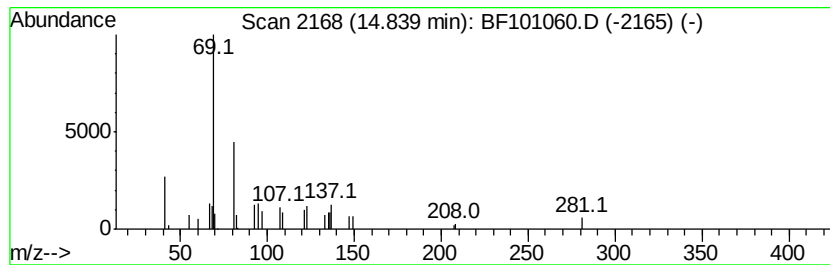
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Peak Number 6 Docosa-2,6,10,14,18-pentaen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	2.01 ng	31460	Perylene-d12	15.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	91
2		Squalene	410	C30H50	000111-02-4	91
3		Supraene	410	C30H50	007683-64-9	86
4		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	003790-71-4	64
5		(.+/-)-Lavandulol, pentafluorop...	300	C13H17F5O2	1000352-63-1	59



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.07	5.07	4.4	ng	63348	1	6.85	289069	20.0
unknown6.62	6.62	89.3	ng	1289990	1	6.85	289069	20.0
Benzoic acid, 2-(...	9.59	2.1	ng	47020	3	9.88	444187	20.0
3-Chloropropionic...	13.84	2.2	ng	47451	5	14.02	437677	20.0
Docosa-2,6,10,14,...	14.84	2.0	ng	31460	6	15.48	313220	20.0