

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100815\
 Data File : BF082131.D
 Acq On : 8 Oct 2015 17:05
 Operator : UM/IZ
 Sample : G3945-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MGP215BR-20-21

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-BF092815.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	5.040	251	254	266	rVB	610167	803470	41.05%	4.576%
2	5.452	288	290	293	rBV	1177012	1482000	75.72%	8.441%
3	6.492	378	381	390	rBV	1502709	1570623	80.25%	8.946%
4	6.629	390	393	395	rBV	1800486	1754038	89.62%	9.991%
5	6.857	411	413	416	rBV	422262	443736	22.67%	2.527%
6	7.017	424	427	429	rBV	1530844	1366138	69.80%	7.781%
7	7.429	460	463	465	rBV	1038092	970284	49.57%	5.526%
8	8.149	523	526	528	rBV	678803	603764	30.85%	3.439%
9	9.098	606	609	611	rBV	18052	21252	1.09%	0.121%
10	9.223	618	620	623	rBV	1867886	1747216	89.27%	9.952%
11	9.623	653	655	657	rBV	262683	248525	12.70%	1.416%
12	9.898	677	679	682	rBV	539517	722287	36.90%	4.114%
13	10.698	746	749	751	rBV	1017747	1244340	63.58%	7.087%
14	10.801	757	758	765	rVB	43943	54725	2.80%	0.312%
15	11.246	795	797	799	rBV	29696	34357	1.76%	0.196%
16	11.395	807	810	812	rBV	485202	703976	35.97%	4.010%
17	11.681	833	835	837	rBV	19468	21039	1.07%	0.120%
18	11.921	854	856	858	rBV	17191	26449	1.35%	0.151%
19	12.972	946	948	951	rBV	1443436	1957235	100.00%	11.148%
20	13.864	1024	1026	1037	rBV	205441	272798	13.94%	1.554%
21	14.024	1038	1040	1043	rBV	490159	733006	37.45%	4.175%
22	14.492	1076	1081	1087	rBV3	12387	31956	1.63%	0.182%
23	14.778	1104	1106	1110	rBV	86996	104157	5.32%	0.593%
24	15.475	1164	1167	1176	rBV	401020	608049	31.07%	3.463%
25	15.967	1207	1210	1214	rBV	11819	31557	1.61%	0.180%

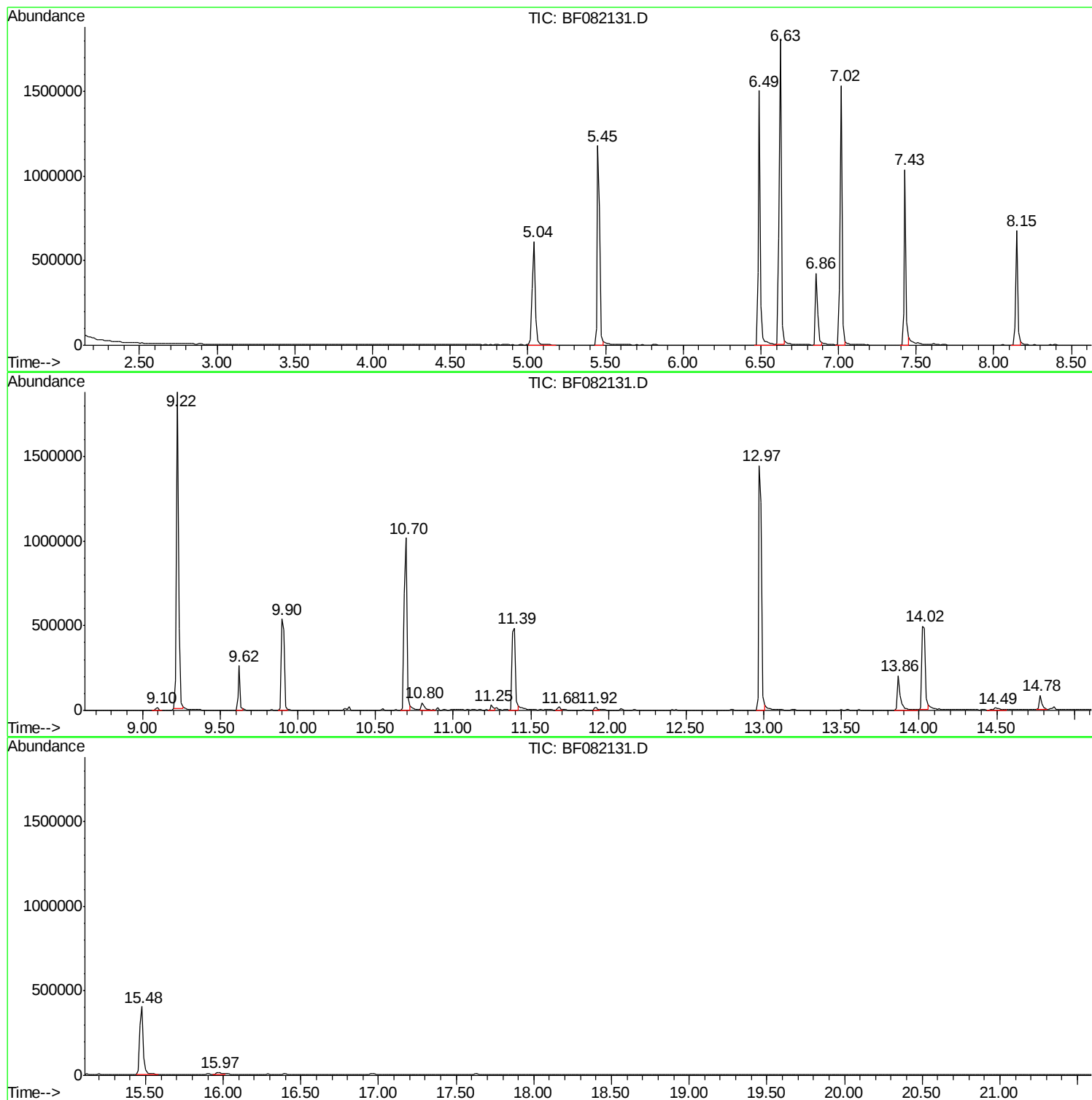
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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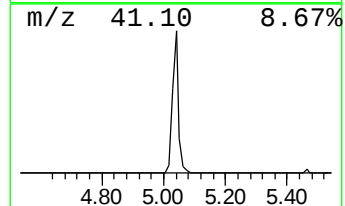
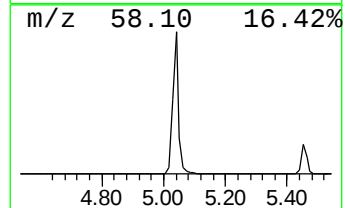
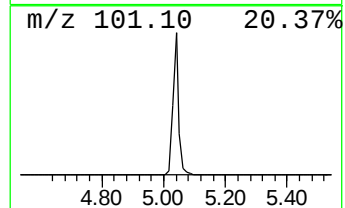
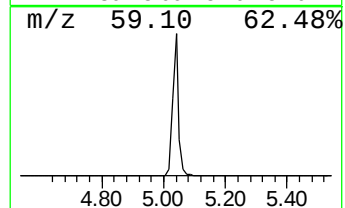
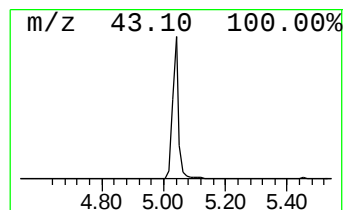
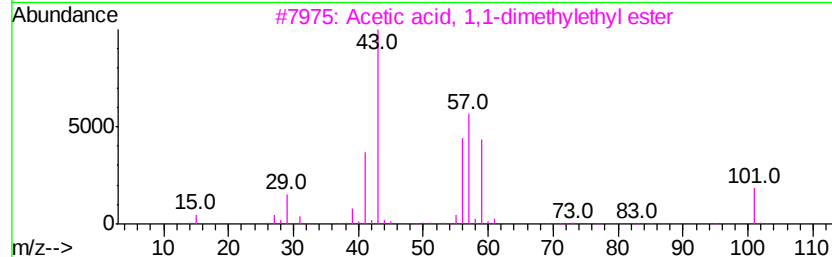
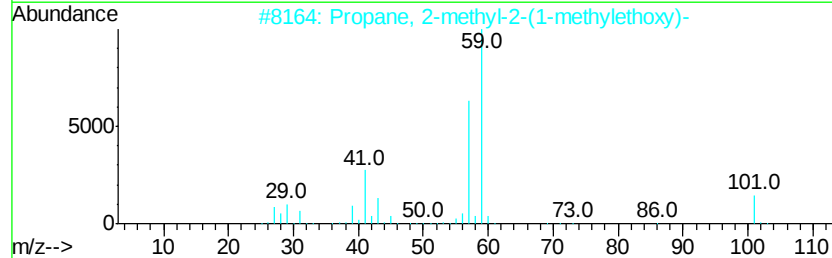
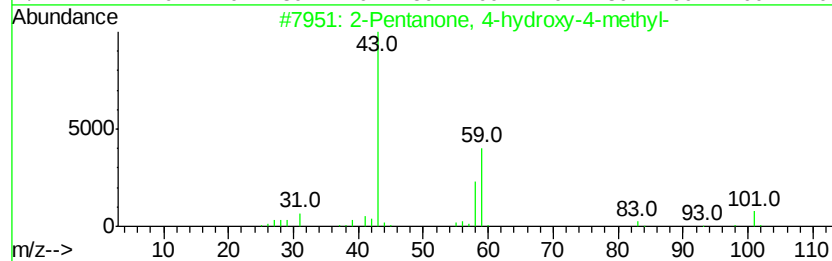
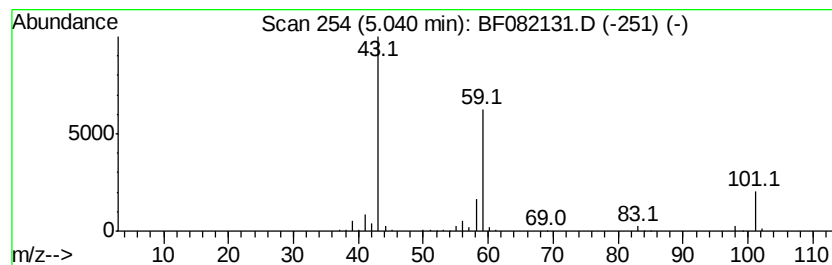
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	36.21 ng	803470	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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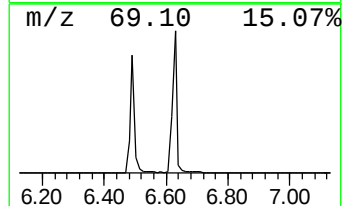
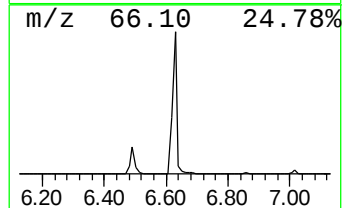
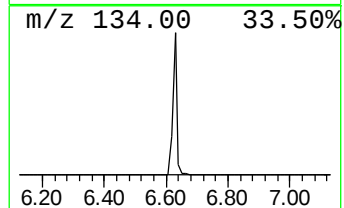
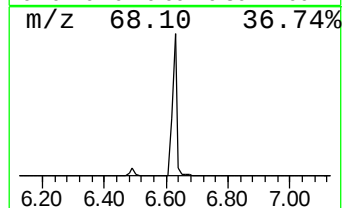
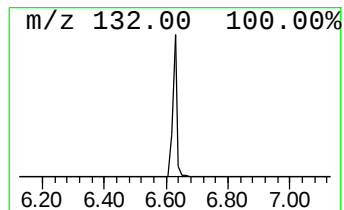
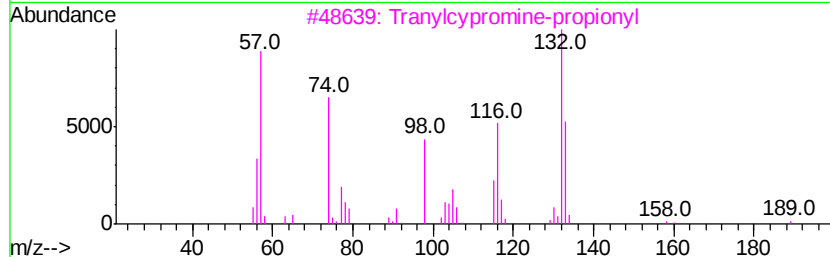
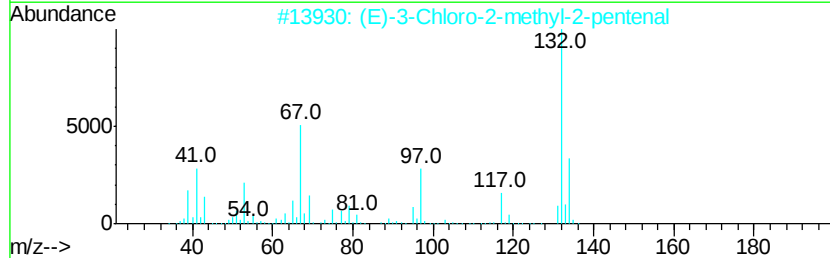
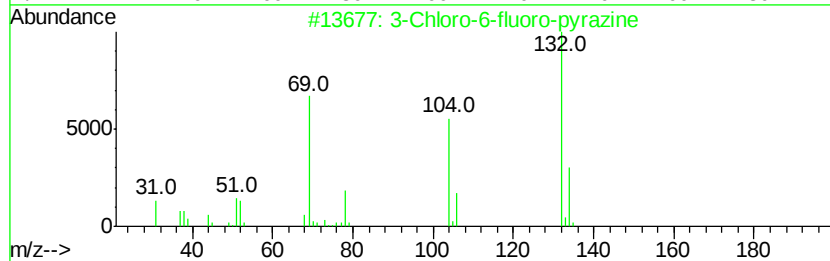
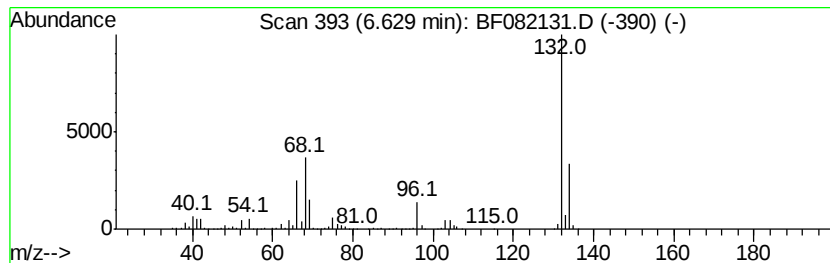
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	79.06 ng	1754040	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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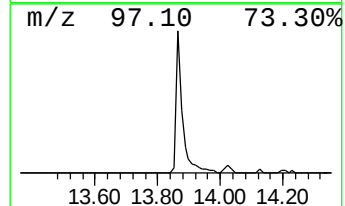
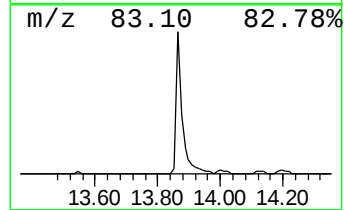
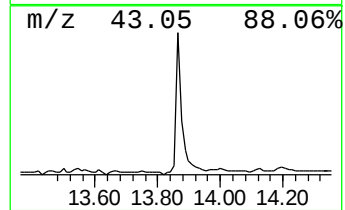
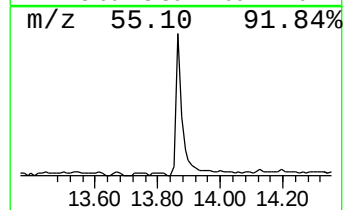
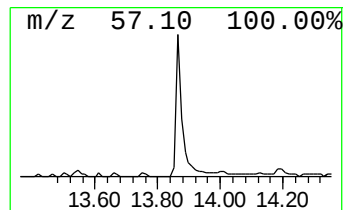
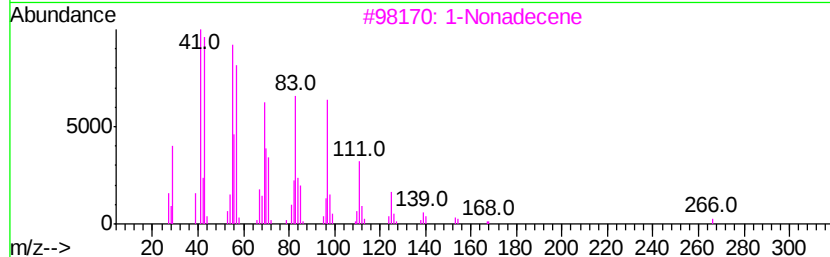
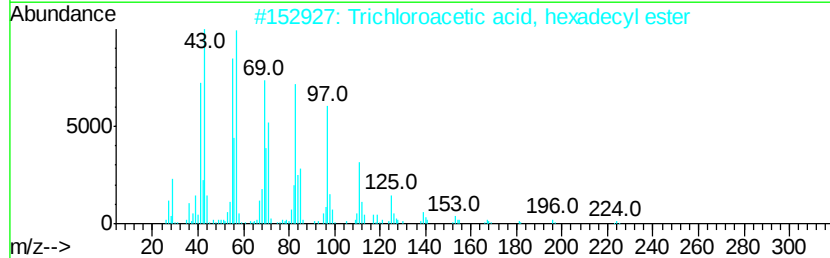
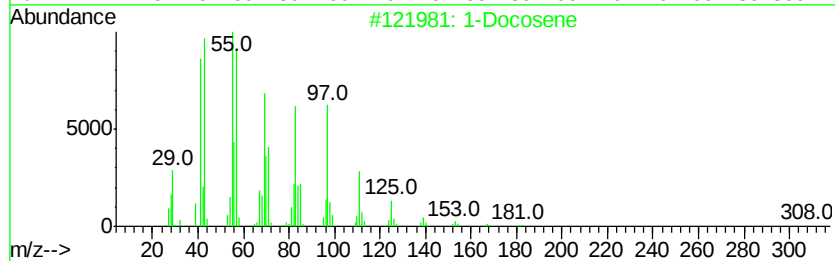
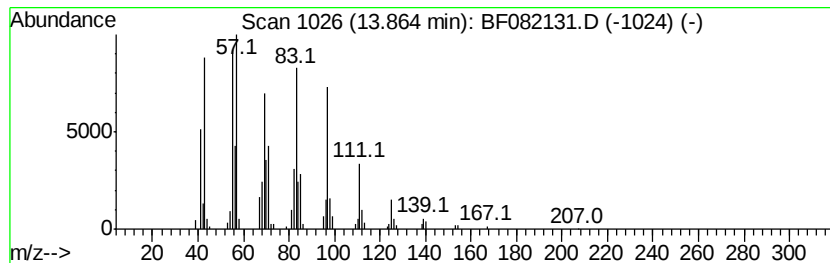
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Peak Number 4 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	7.44 ng	272798	Chrysene-d12	14.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			1-Nonadecene	266	C19H38	018435-45-5	91
4			3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91



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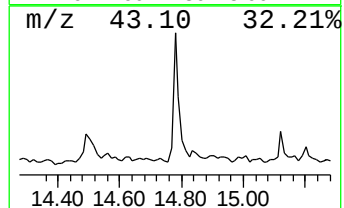
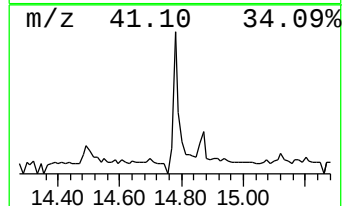
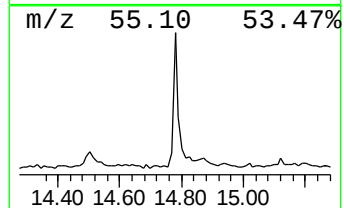
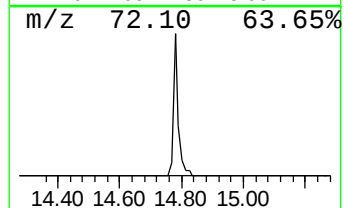
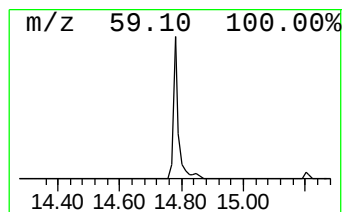
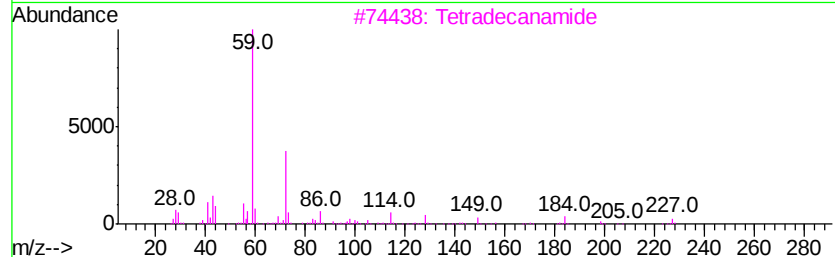
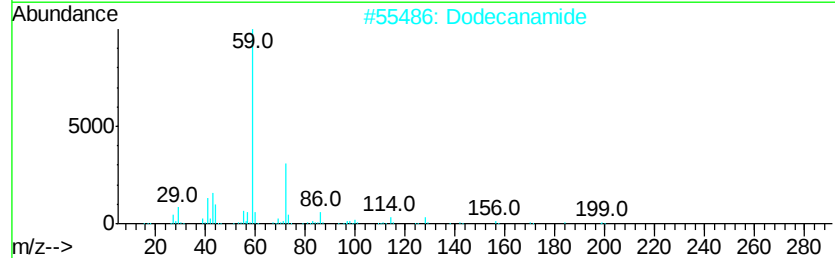
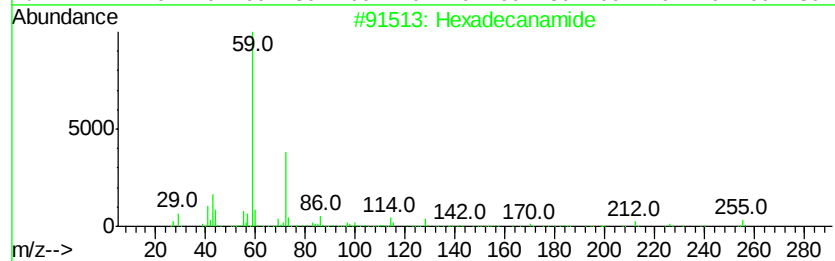
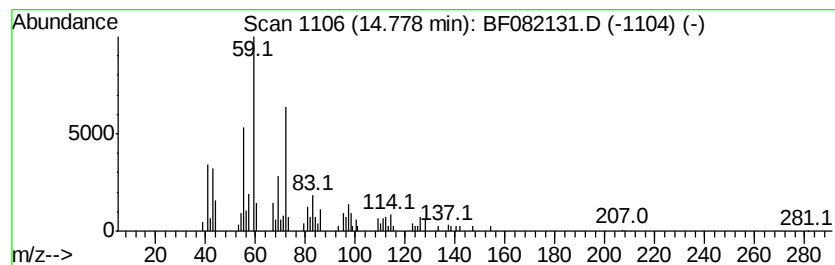
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Peak Number 5 Hexadecanamide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.78	3.43 ng	104157	Perylene-d12	15.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanamide		255	C16H33NO	000629-54-9	64
2	Dodecanamide		199	C12H25NO	001120-16-7	59
3	Tetradecanamide		227	C14H29NO	000638-58-4	53
4	Acetic acid, cyanohydroxvimino-,...		128	C4H4N2O3	061295-92-9	46
5	Octanamide		143	C8H17NO	000629-01-6	45



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.04	36.2	ng	803470	1	6.86	443736	20.0
unknown6.63	6.63	79.1	ng	1754040	1	6.86	443736	20.0
1-Docosene	13.86	7.4	ng	272798	5	14.04	733006	20.0
Hexadecanamide	14.78	3.4	ng	104157	6	15.48	608049	20.0