

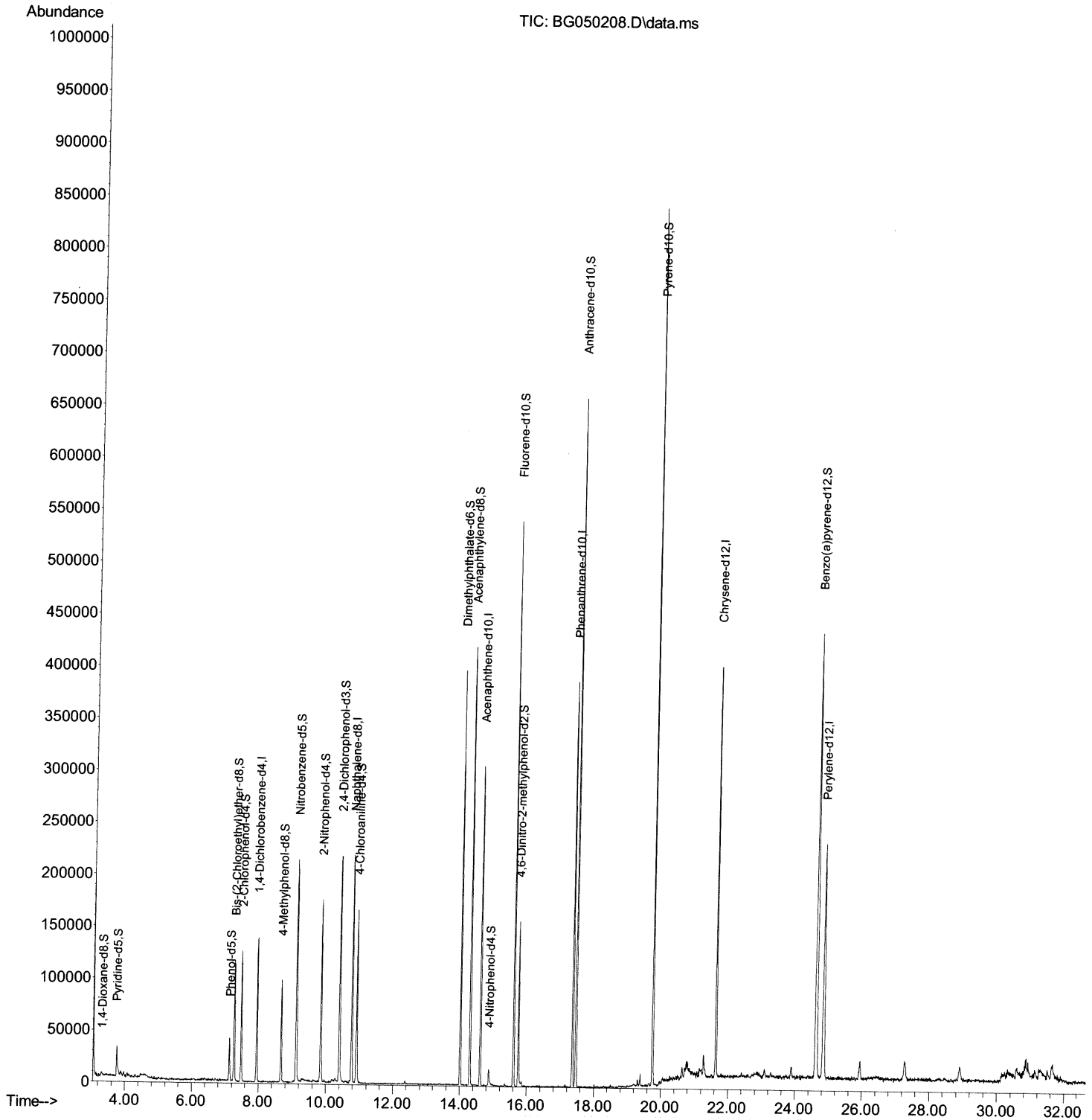
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050208.D
Acq On : 8 Sep 2021 21:27
Operator : CG/JU
Sample : M3651-15
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKM9

Manual Integrations
APPROVED

mohammad
9/9/2021 11:46:27 AM

Quant Time: Sep 09 00:57:03 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
Qlast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

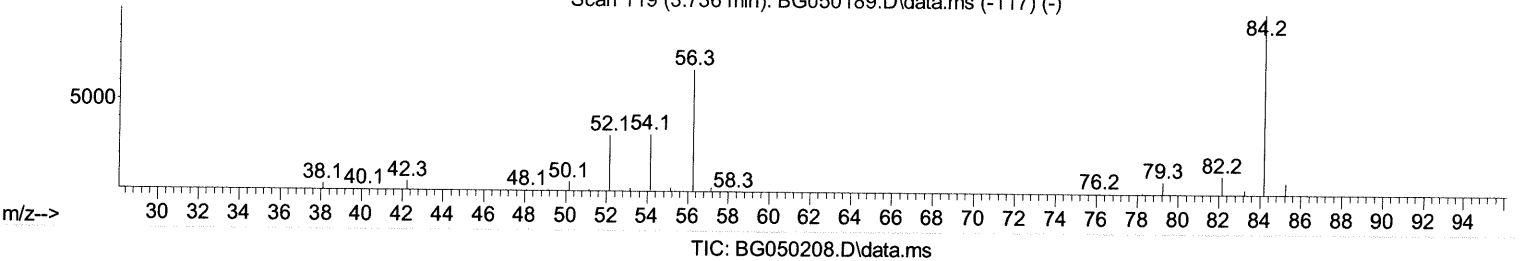
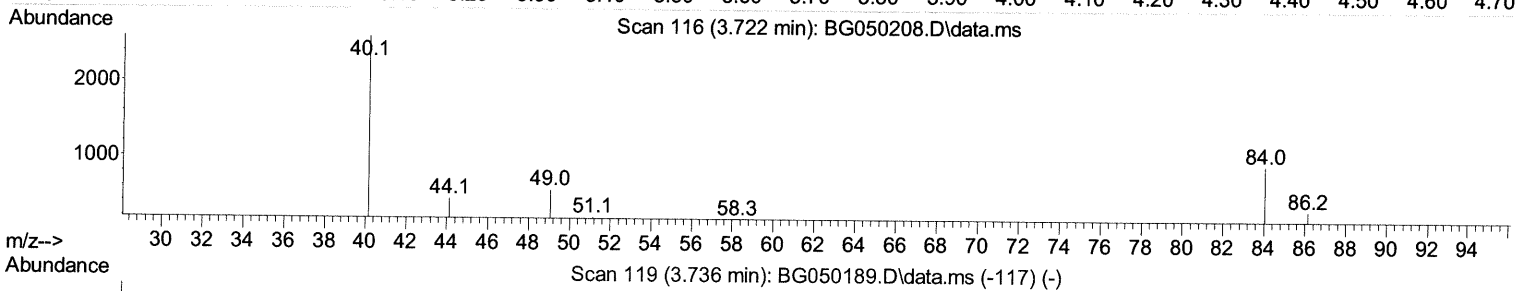
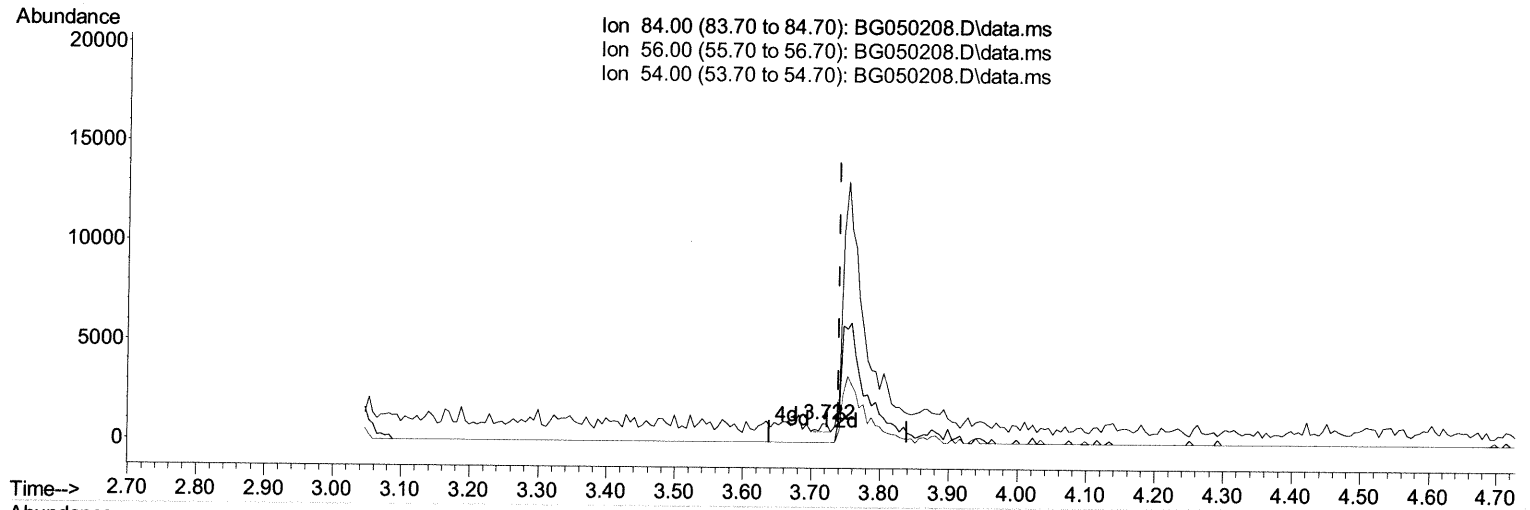
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(4) Pyridine-d5 (S)

3.722min (-0.016) 0.12 ng/ul

response 359

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	0.00#
54.00	37.00	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

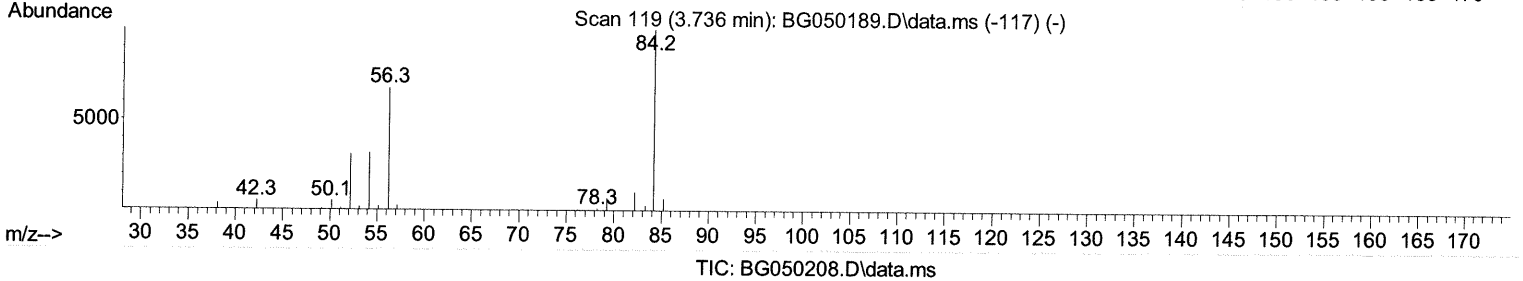
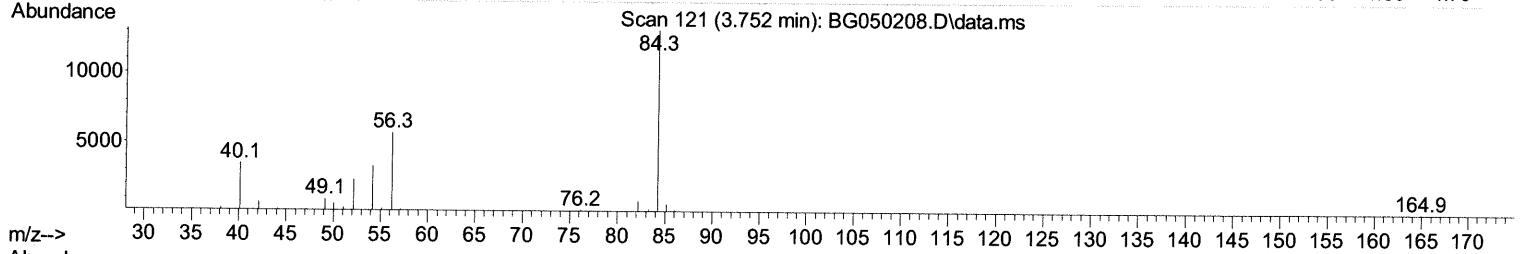
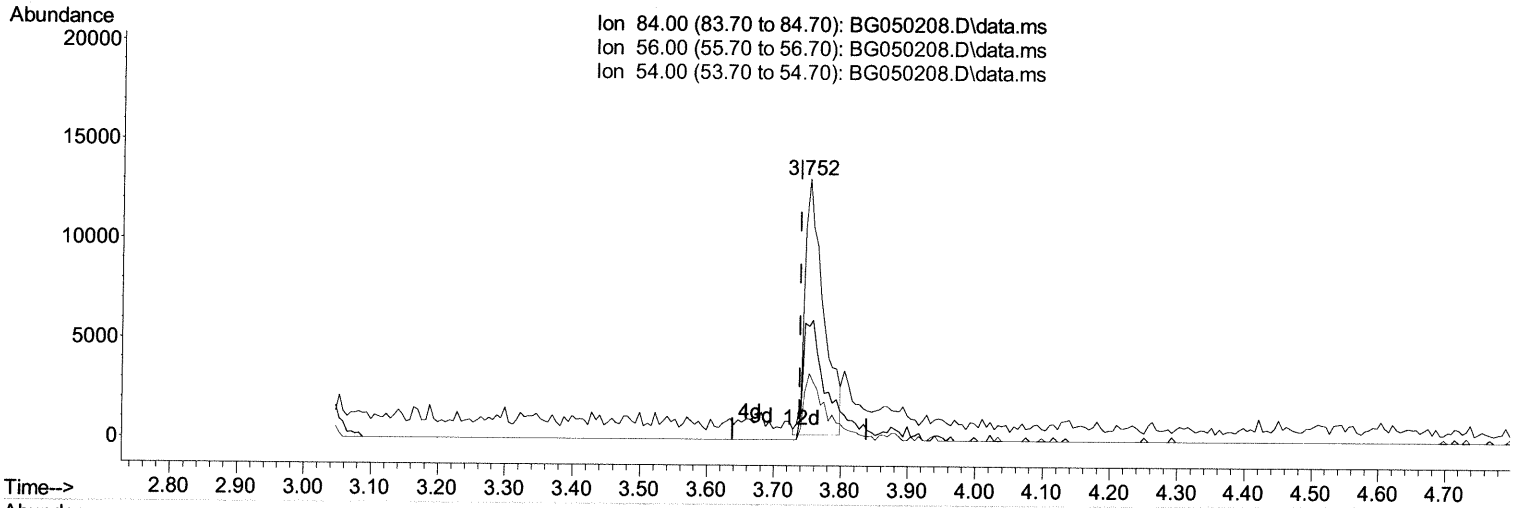
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TIC: BG050208.D\data.ms

(4) Pyridine-d5 (S)

3.752min (+ 0.013) 8.64 ng/ul m 09/10/21 34

response 24997

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	43.52
54.00	37.00	25.34#
0.00	0.00	0.00

Quantitation Report (Qedit)

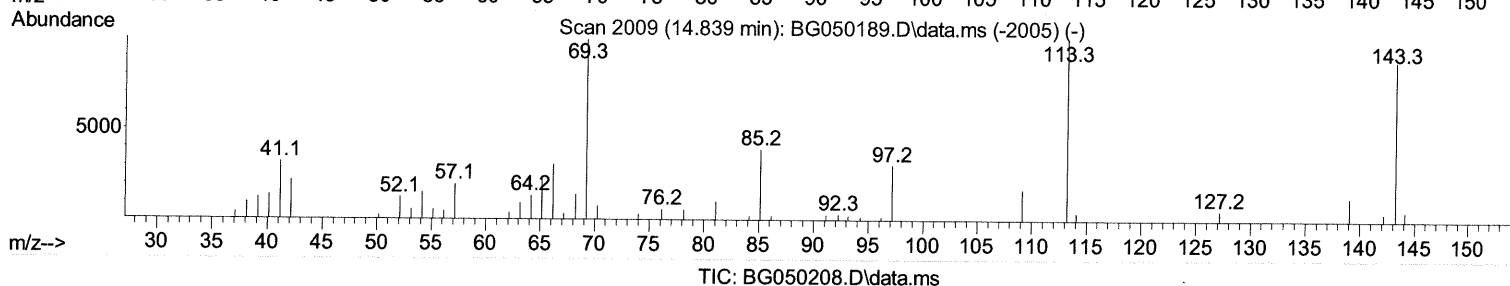
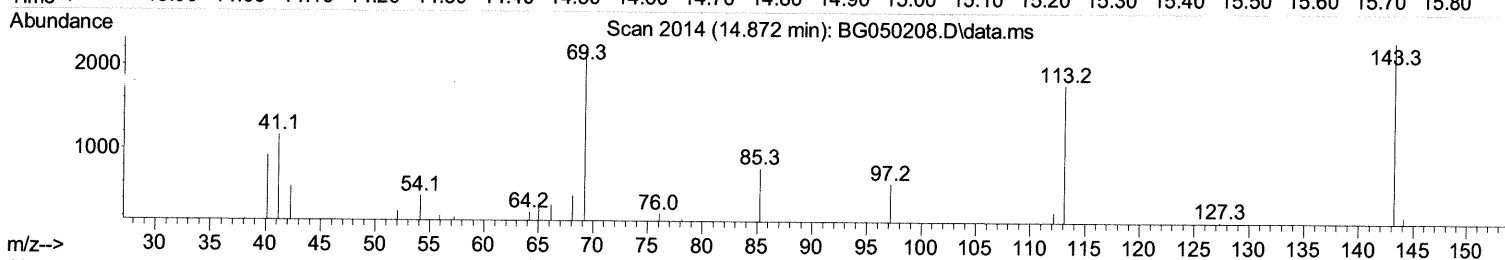
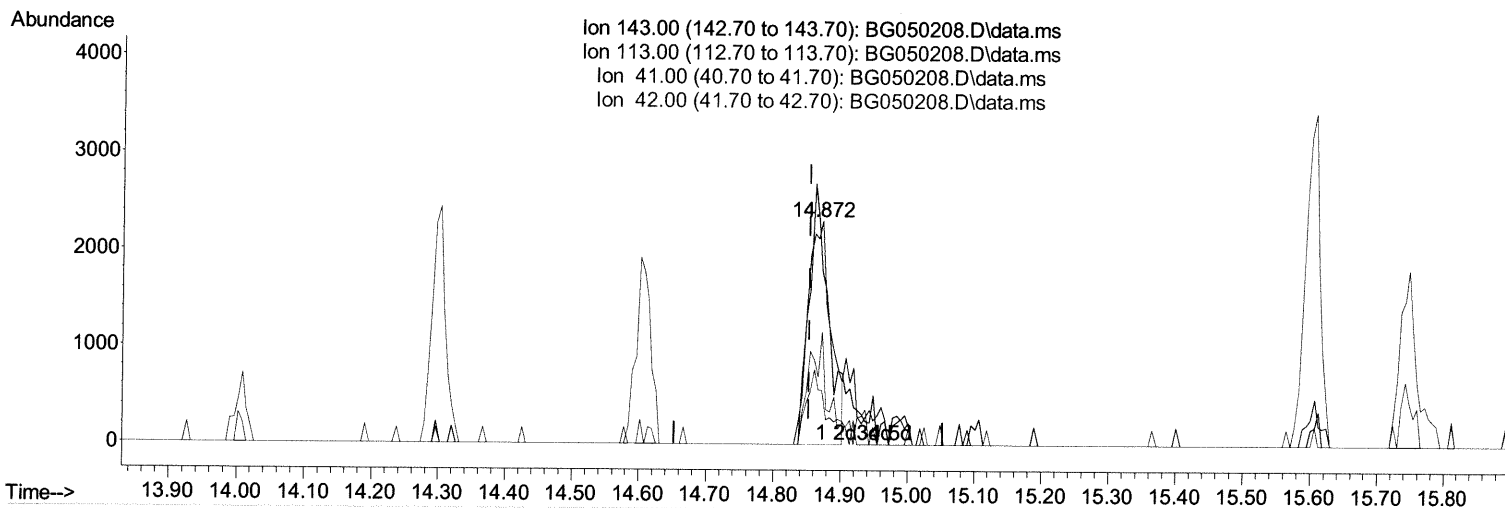
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(54) 4-Nitrophenol-d4 (S)

14.872min (+ 0.019) 5.55 ng/ul

response 5517

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	121.00	77.54#
41.00	45.30	50.65
42.00	25.60	24.15

Quantitation Report (Qedit)

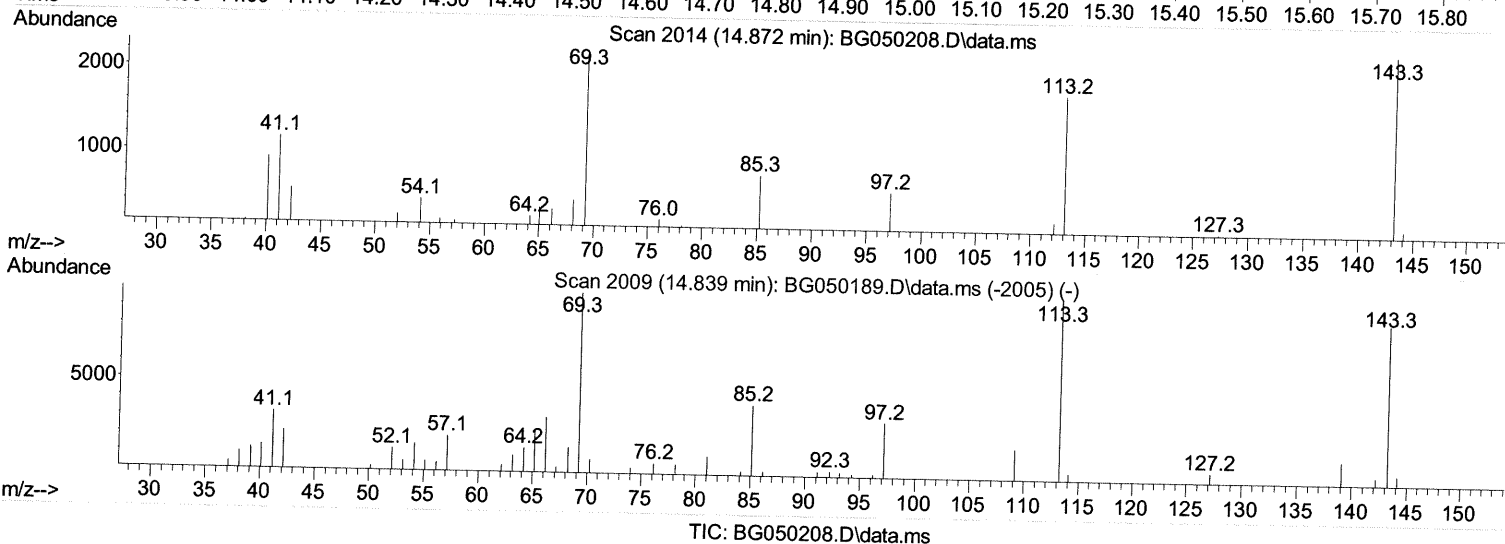
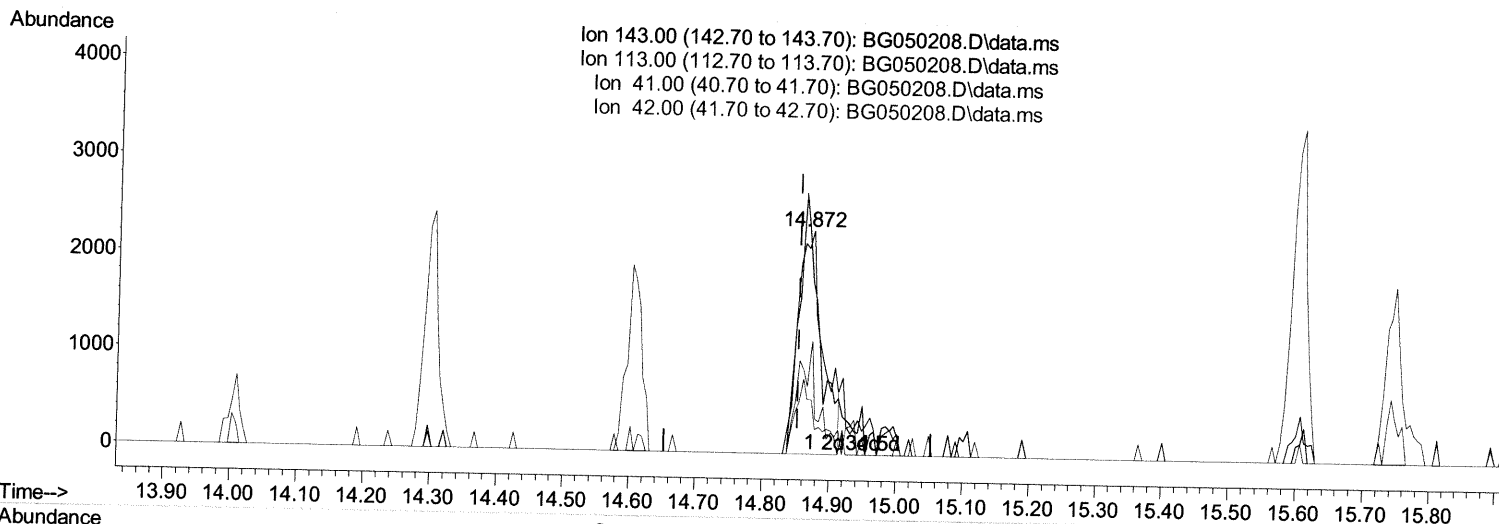
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	40289	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.766	136	198584	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.608	164	109109	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.363	188	239064	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	243528	20.000	ng/ul	0.00
88) Perylene-d12	24.847	264	246271	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	2240	2.749	ng/uL	0.00
4) Pyridine-d5	3.752	84	24997m	8.641	ng/ul	0.01
7) Phenol-d5	7.129	99	27951	8.935	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.265	67	57757	31.615	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.482	132	59098	22.764	ng/ul	0.00
15) 4-Methylphenol-d8	8.674	113	40130	16.316	ng/ul	0.00
21) Nitrobenzene-d5	9.121	128	52697	38.661	ng/ul	0.00
24) 2-Nitrophenol-d4	9.843	143	55091	33.239	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.402	165	87350	27.613	ng/ul	0.00
31) 4-Chloroaniline-d4	10.913	131	94671	24.291	ng/ul	0.00
46) Dimethylphthalate-d6	14.008	166	257163	31.214	ng/ul	-0.01
49) Acenaphthylene-d8	14.302	160	308504	31.448	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.872	143	6056m	6.092	ng/ul	0.02
60) Fluorene-d10	15.606	176	226264	33.076	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.741	200	37802	26.214	ng/ul	-0.01
73) Anthracene-d10	17.463	188	387610	36.117	ng/ul	0.00
81) Pyrene-d10	19.754	212	472949	38.989	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.629	264	468035	35.901	ng/ul	-0.01

Target Compounds

Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed