

Quantitation Report (QT Reviewed)

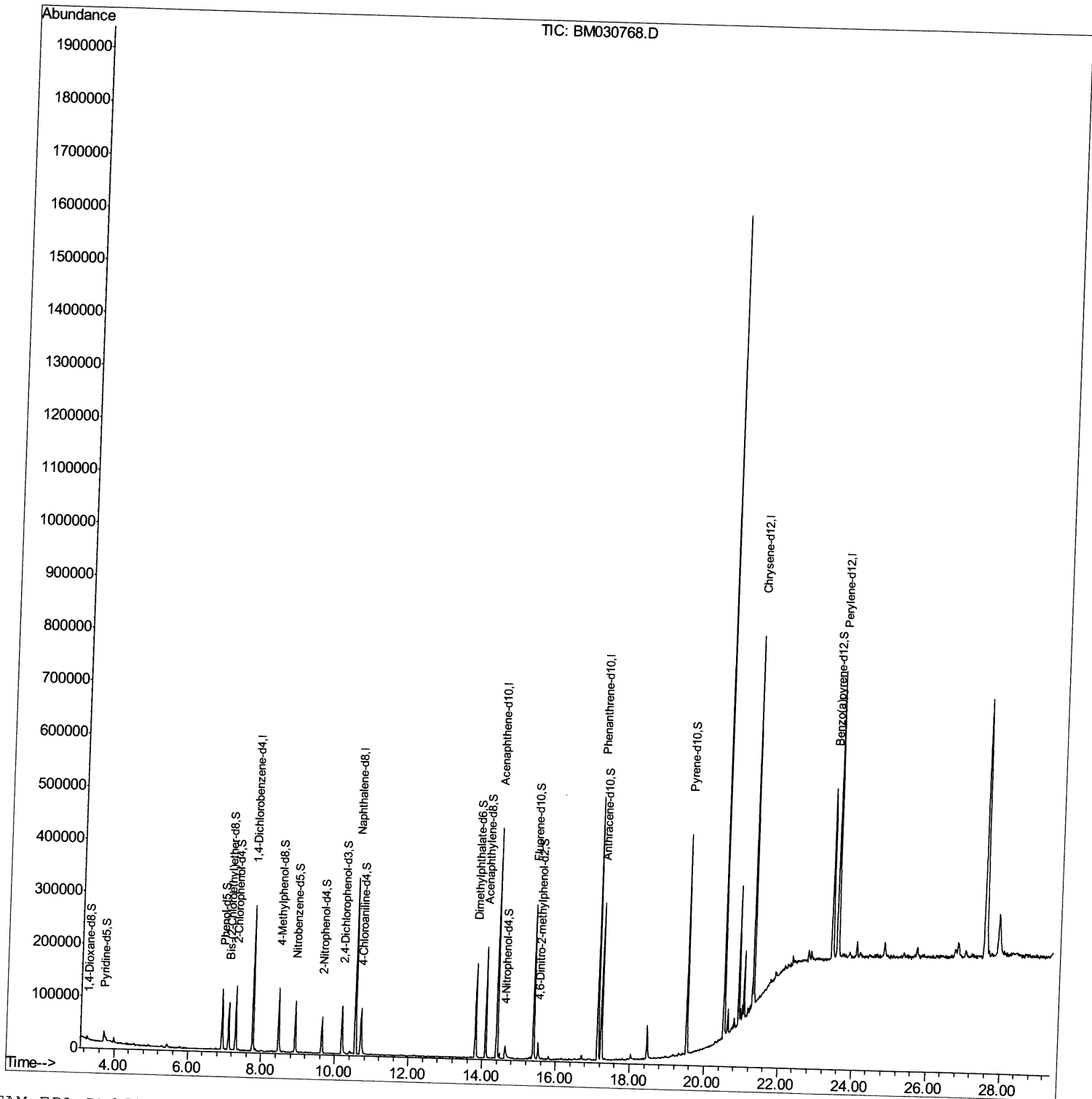
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM063021\
 Data File : BM030768.D
 Acq On : 02 Jul 2021 06:52
 Operator : CG/JU
 Sample : M2825-20
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGDR9

Manual Integrations
 APPROVED

mohammad
 7/2/2021 12:04:25 PM

Quant Time: Jul 02 07:40:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 02 04:13:39 2021
 Response via : Initial Calibration



Quantitation Report (Qedit)

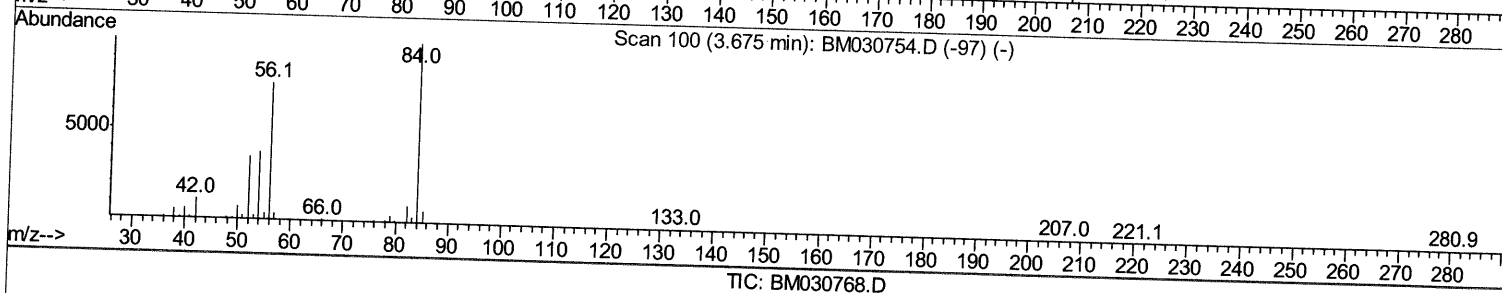
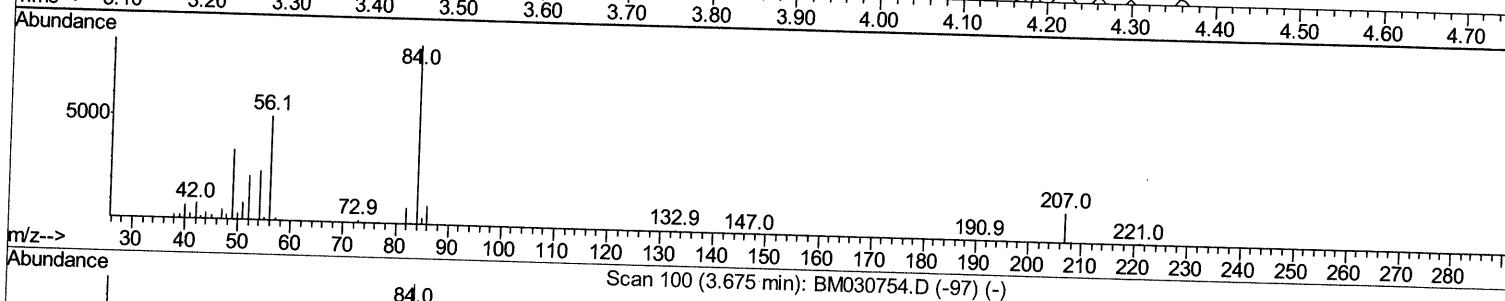
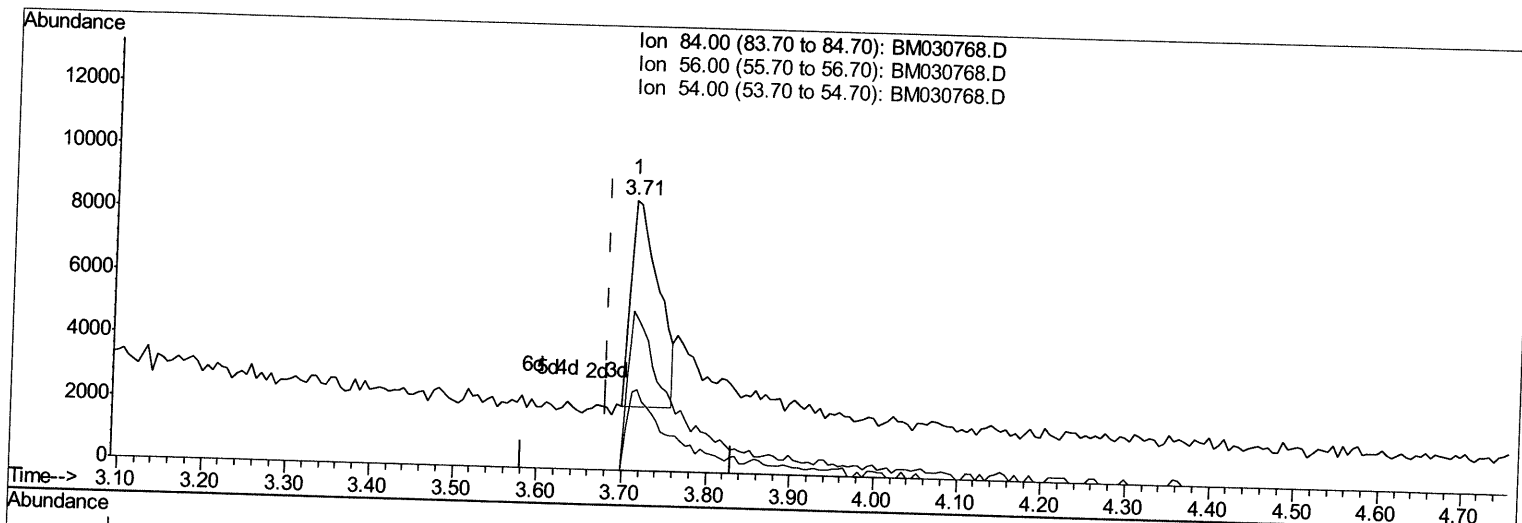
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(4) Pyridine-d5 (S)

3.710min (+0.029) 3.20ng/ul

response 15082

Ion	Exp%	Act%
84.00	100	100
56.00	73.70	58.99
54.00	35.80	29.00
0.00	0.00	0.00

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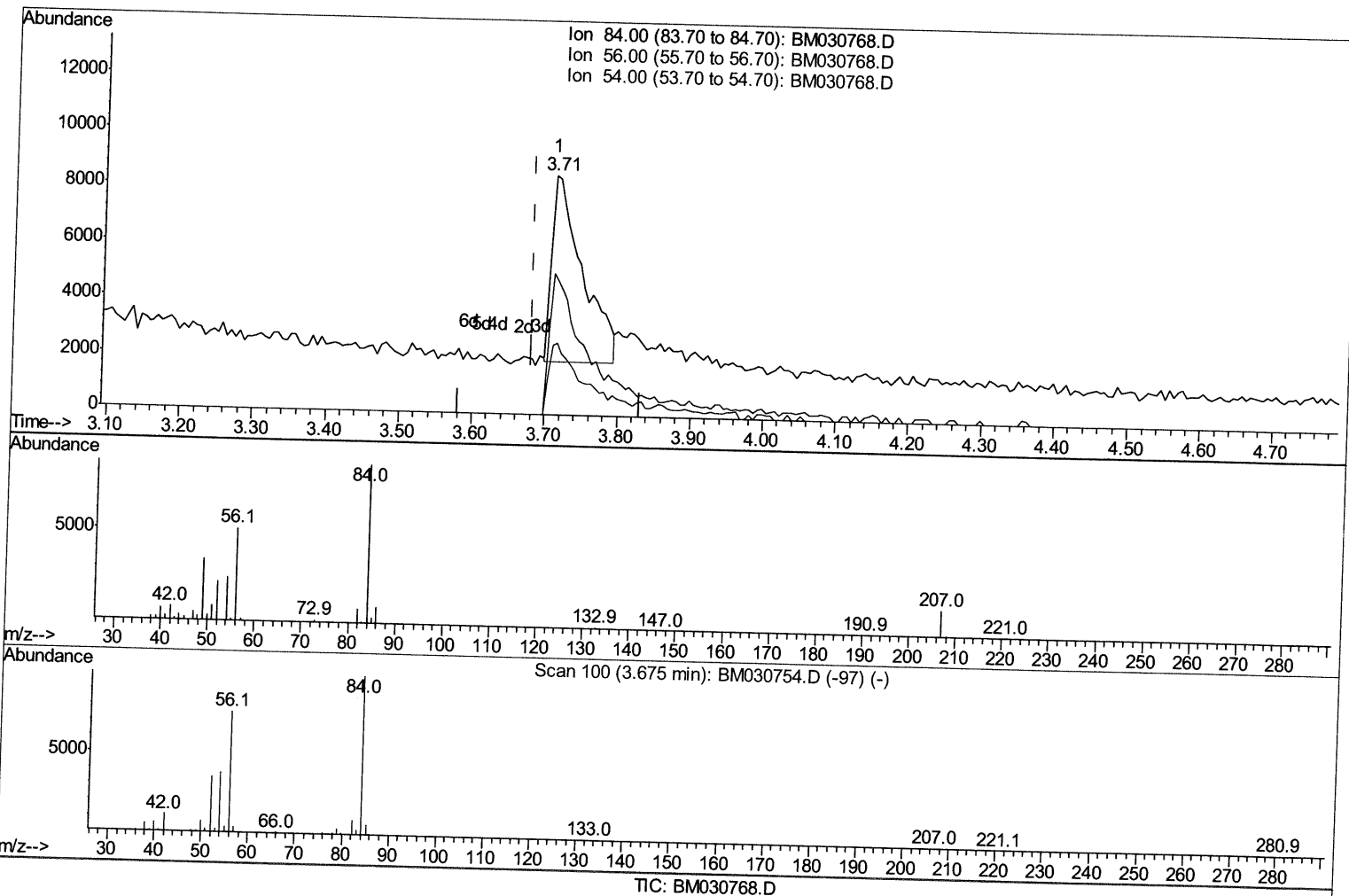
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(4) Pyridine-d5 (S)

3.710min (+0.029) 4.06ng/ul m

response 19128

Ion	Exp%	Act%
84.00	100	100
56.00	73.70	58.99
54.00	35.80	29.00
0.00	0.00	0.00

347/7/21

Quantitation Report (Qedit)

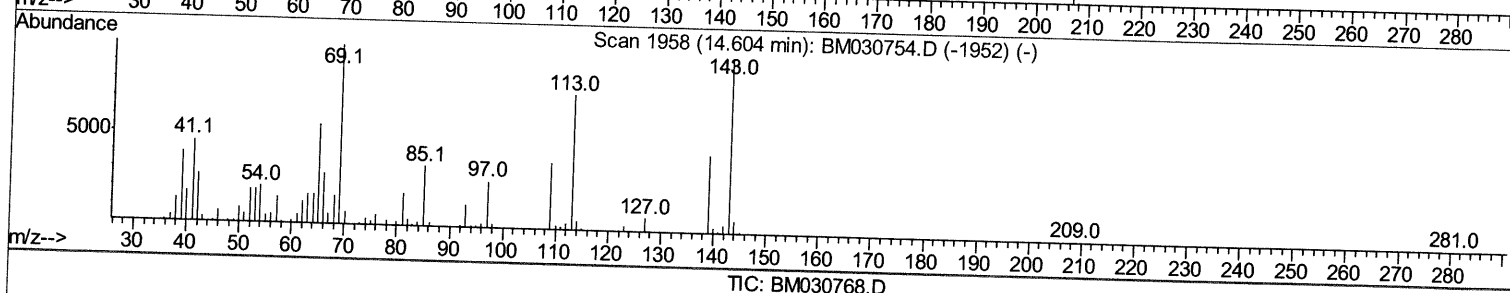
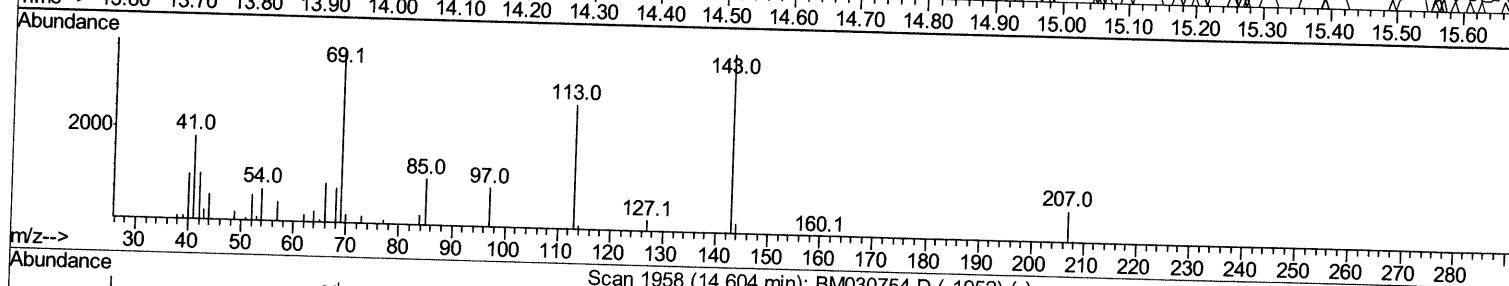
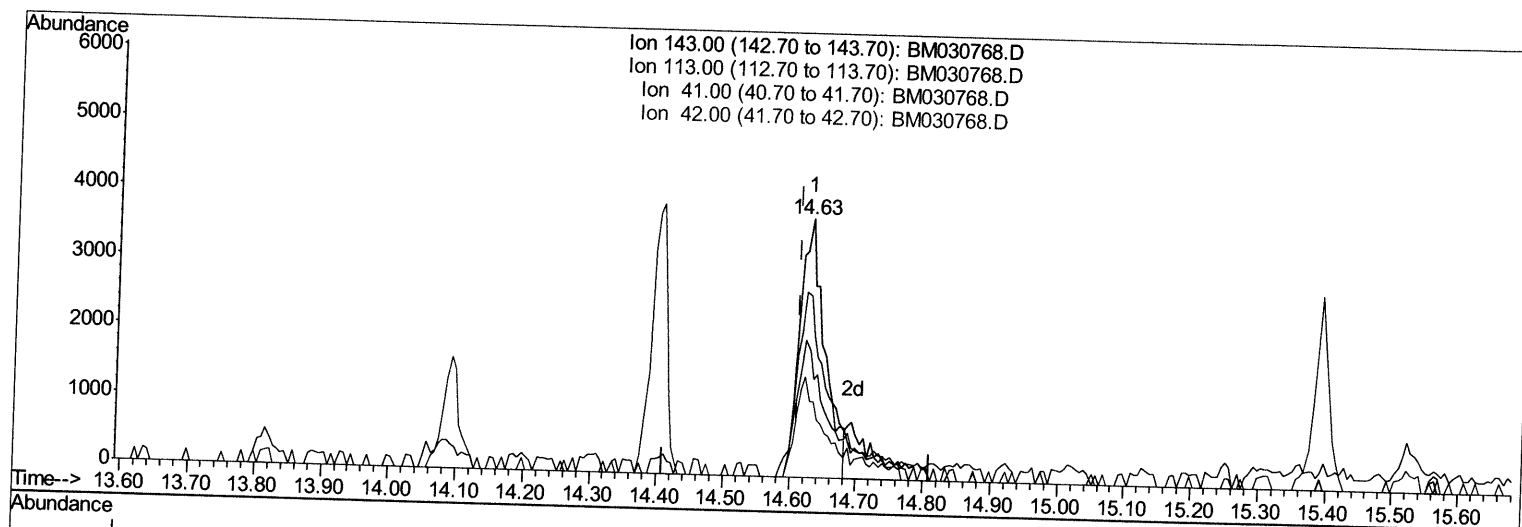
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(54) 4-Nitrophenol-d4 (S)

14.627min (+0.018) 5.02ng/ul

response 9873

Ion	Exp%	Act%
143.00	100	100
113.00	74.50	70.54
41.00	44.10	48.97
42.00	30.20	29.24

Quantitation Report (Qedit)

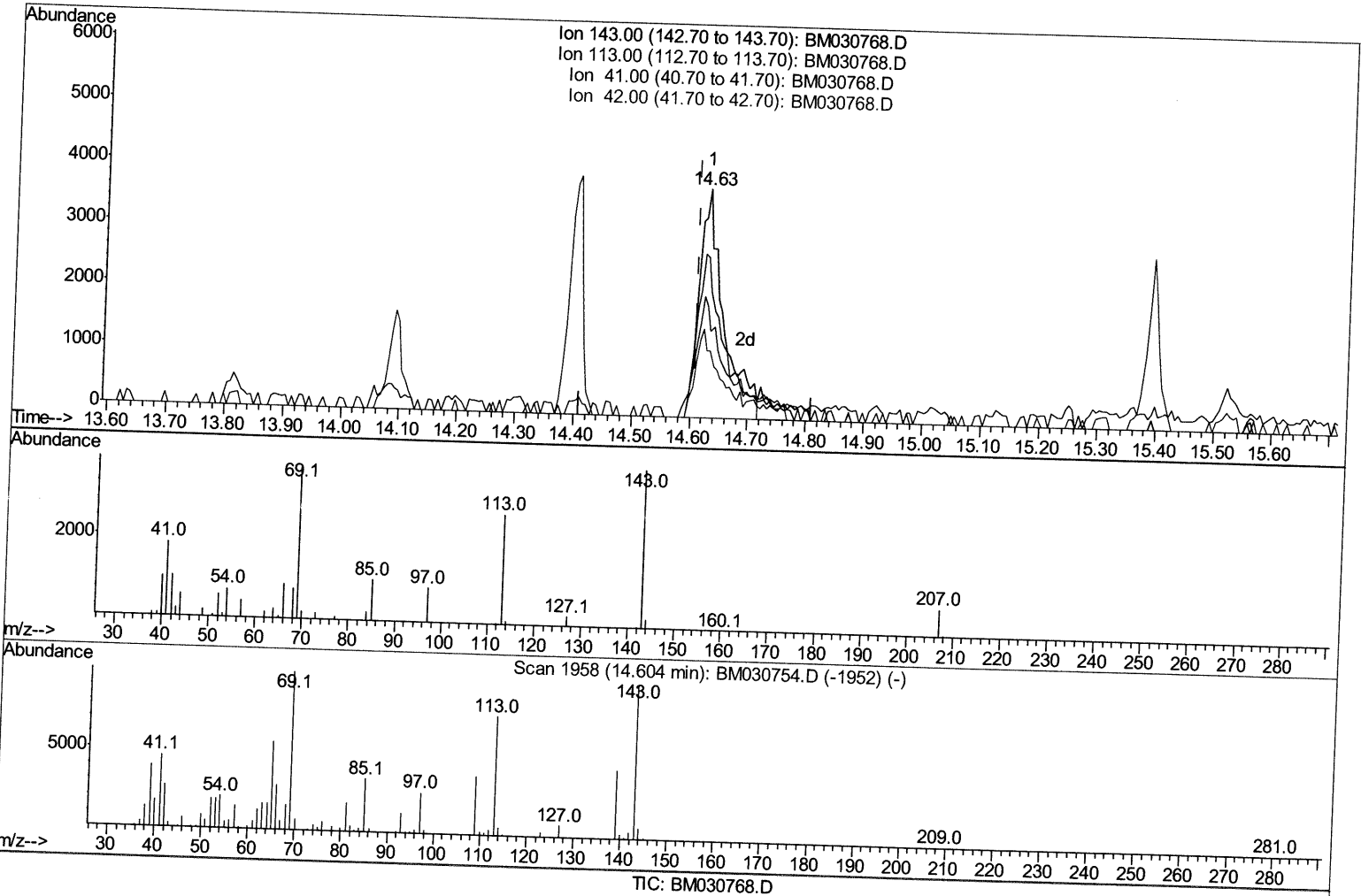
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(54) 4-Nitrophenol-d4 (S)

14.627min (+0.018) 5.67ng/ul m

response 11151

Ion	Exp%	Act%
143.00	100	100
113.00	74.50	70.54
41.00	44.10	48.97
42.00	30.20	29.24

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TIC: BM030768.D

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	71115	20.00	ng/ul	-0.01
20) Naphthalene-d8	10.56	136	269883	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.40	164	150231	20.00	ng/ul	-0.01
64) Phenanthrene-d10	17.14	188	294878	20.00	ng/ul	0.00
79) Chrysene-d12	21.31	240	322947	20.00	ng/ul	0.00
88) Perylene-d12	23.56	264	371293	20.00	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.27	96	4301	2.45	ng/uL	0.00
4) Povidine-d5	3.71	84	19128m	4.06	ng/ul	0.03
7) Phenol-d5	6.94	99	61646	10.05	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.11	67	43168	11.20	ng/ul	0.00
11) 2-Chlorophenol-d4	7.30	132	52872	11.17	ng/ul	0.00
15) 4-Methylphenol-d8	8.47	113	46278	9.70	ng/ul	0.00
21) Nitrobenzene-d5	8.93	128	21323	10.59	ng/ul	0.00
24) 2-Nitrophenol-d4	9.65	143	20798	9.29	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.19	165	39995	9.33	ng/ul	0.00
31) 4-Chloroaniline-d4	10.70	131	55369	9.26	ng/ul	0.00
46) Dimethylphthalate-d6	13.82	166	128729	12.41	ng/ul	0.00
49) Acenaphthylene-d8	14.09	160	151095	11.22	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.63	143	11151m	5.67	ng/ul	0.02
60) Fluorene-d10	15.39	176	114330	12.43	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.52	200	8594	4.45	ng/ul	0.00
73) Anthracene-d10	17.24	188	172992	12.57	ng/ul	0.00
81) Pyrene-d10	19.53	212	201504	11.91	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.41	264	234224	12.13	ng/ul	-0.02

7/2/21

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Target Compounds

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