

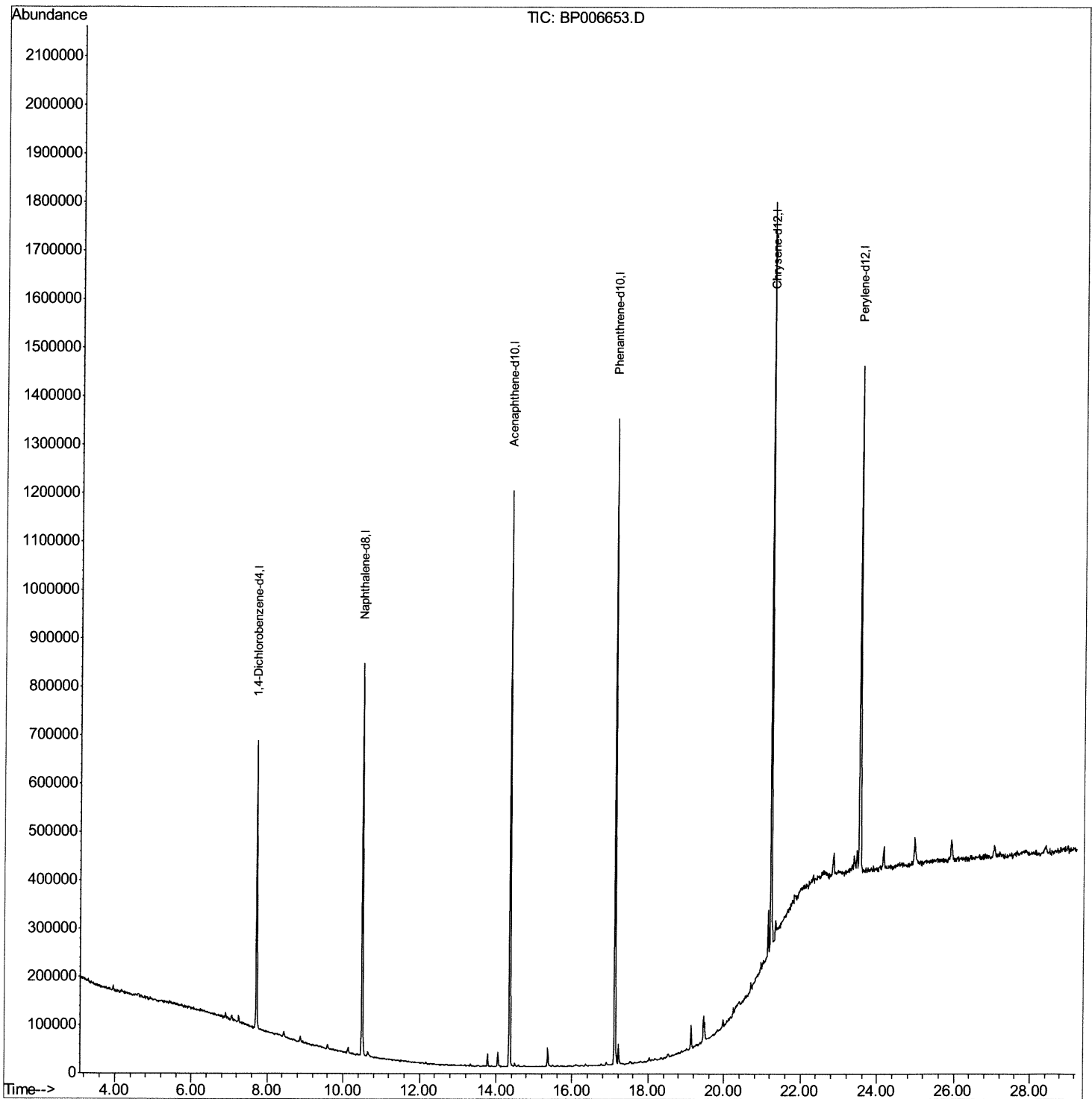
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP081021\
Data File : BP006653.D
Acq On : 11 Aug 2021 13:19
Operator : CG/JU
Sample : M3246-06 30X
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB0

Manual Integrations
APPROVED

mohammad
8/11/2021 9:29:42 PM

Quant Time: Aug 11 14:04:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP081021.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 10 15:09:30 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

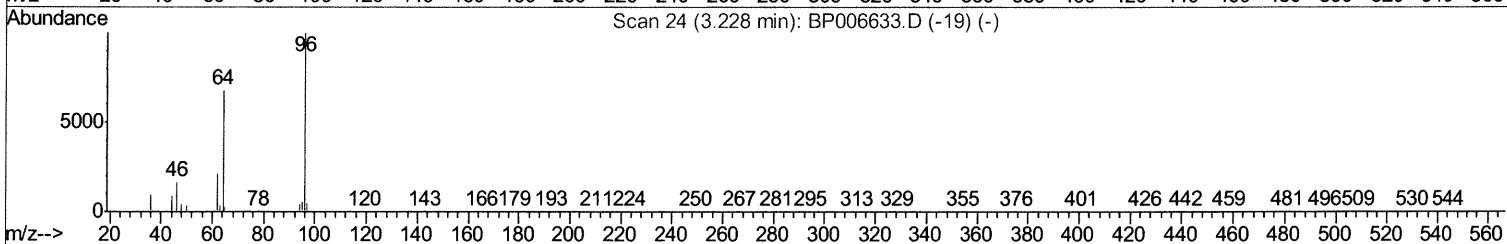
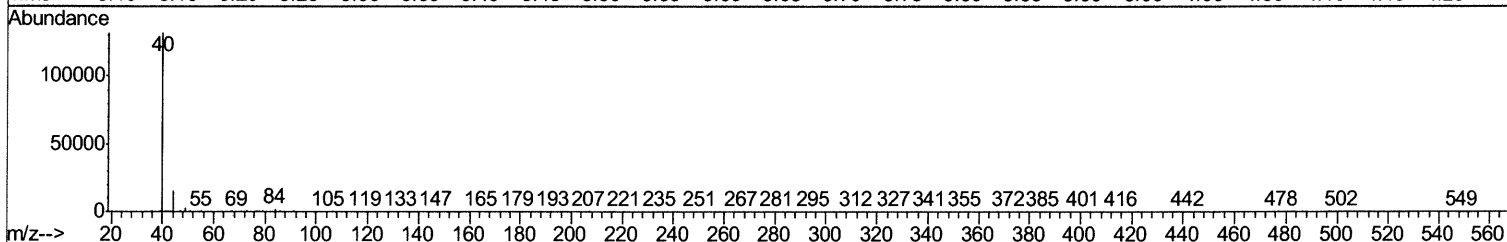
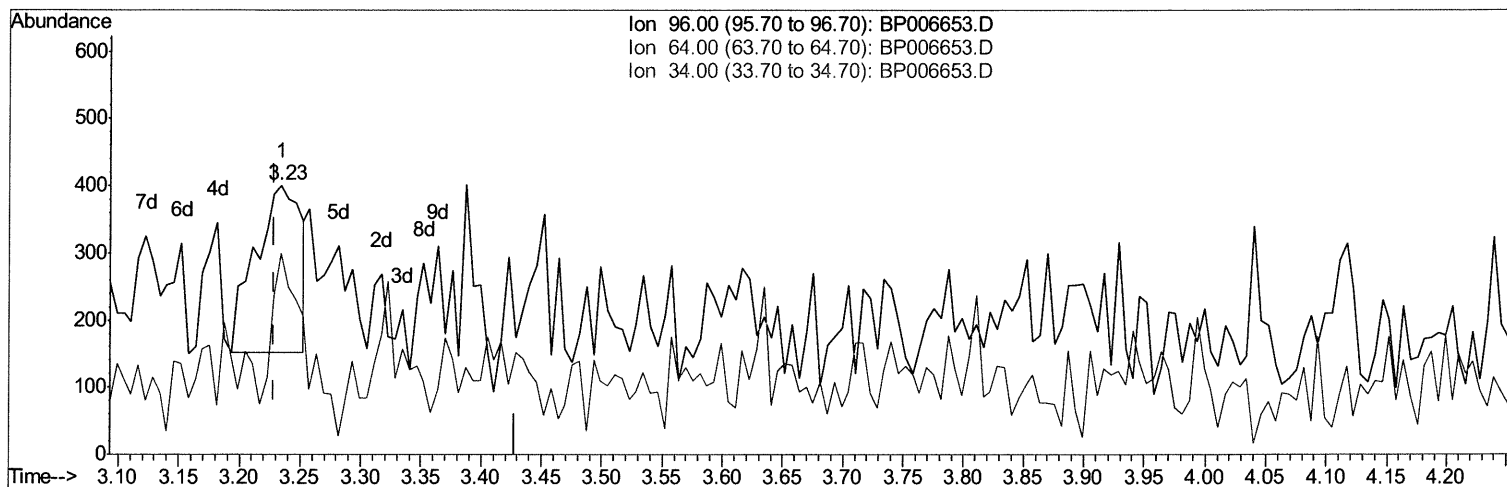
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TIC: BP006653.D

(3) 1,4-Dioxane-d8 (S)

3.234min (+0.006) 0.15ng/uL

response 640

Ion	Exp%	Act%
96.00	100	100
64.00	70.10	74.81
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

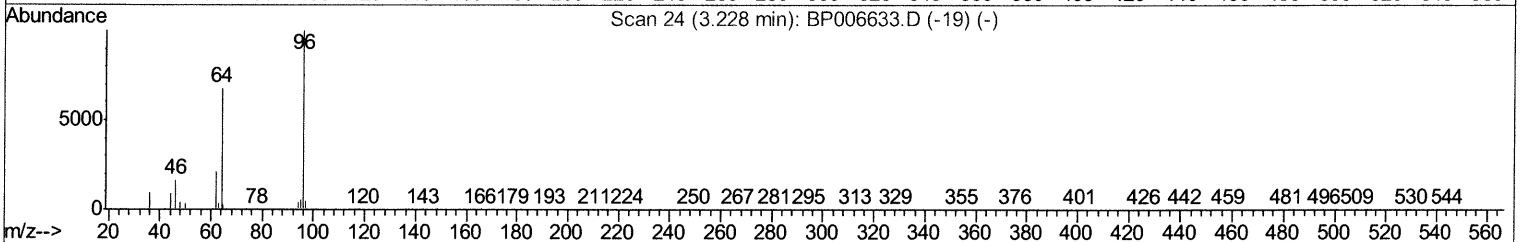
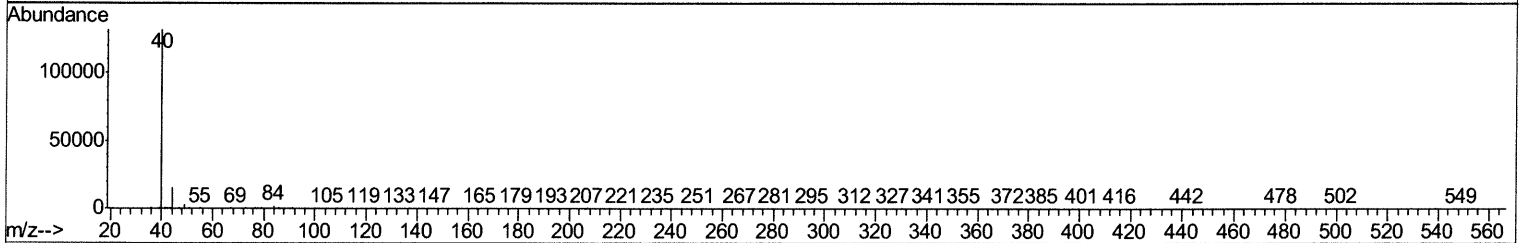
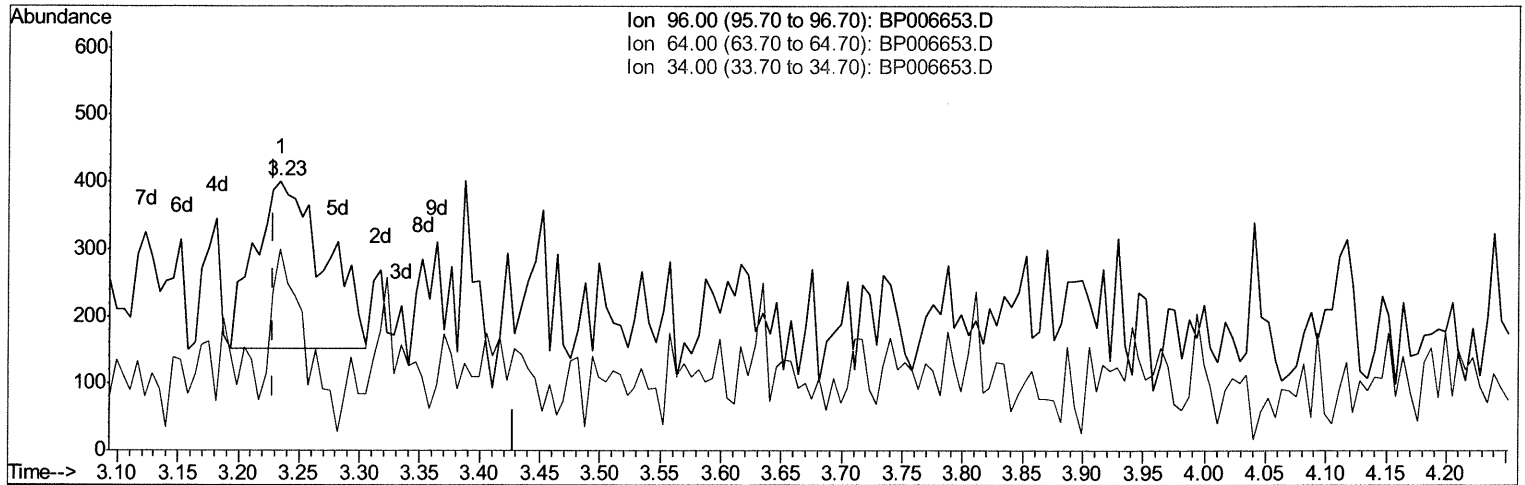
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TIC: BP006653.D

(3) 1,4-Dioxane-d8 (S)

3.234min (+0.006) 0.23ng/uL m

response 998

Ion	Exp%	Act%
96.00	100	100
64.00	70.10	74.81
34.00	0.00	0.00
0.00	0.00	0.00

Ju 8/13/21

Quantitation Report (Qedit)

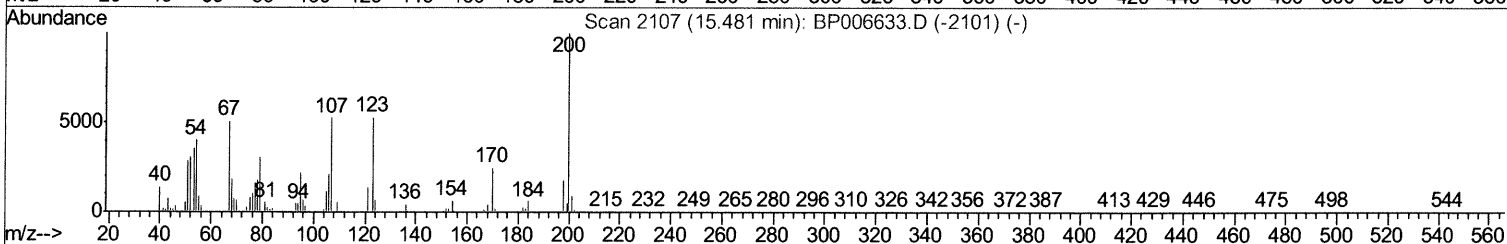
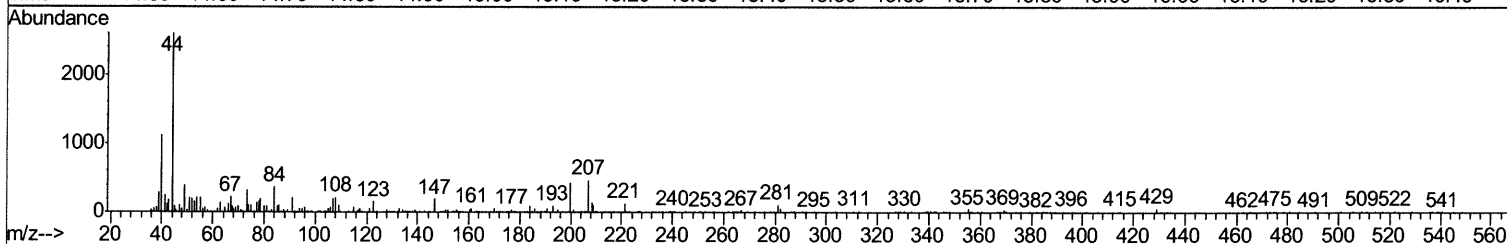
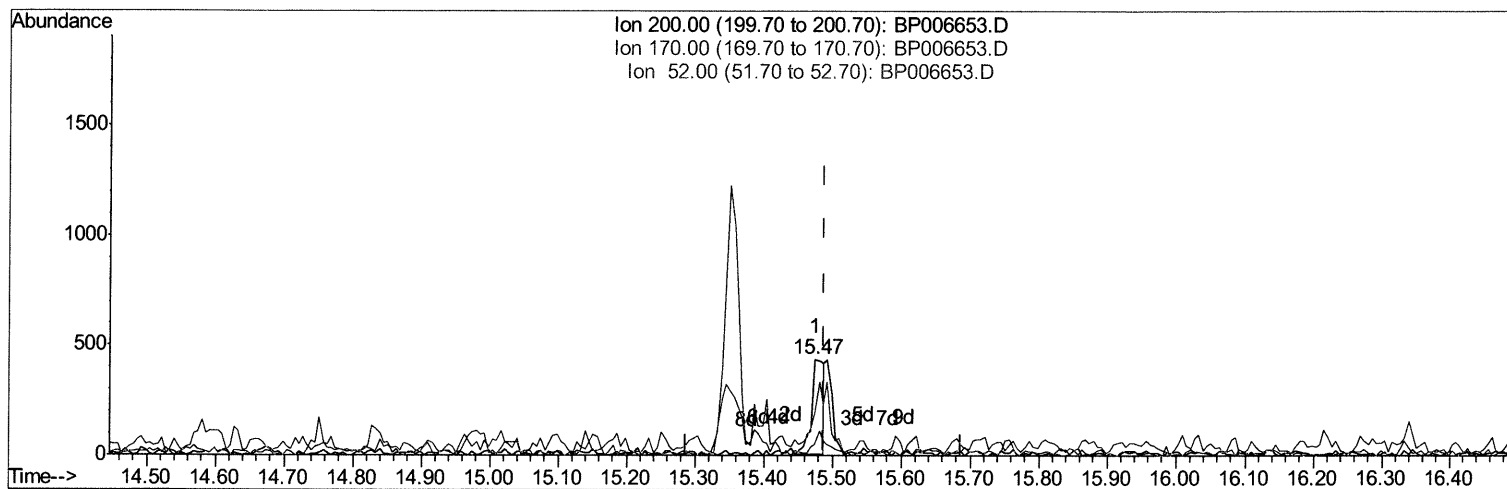
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TIC: BP006653.D

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.475min (-0.012) 0.12ng/ul

response 531

Ion	Exp%	Act%
200.00	100	100
170.00	22.40	11.67#
52.00	47.50	43.71
0.00	0.00	0.00

Quantitation Report (Qedit)

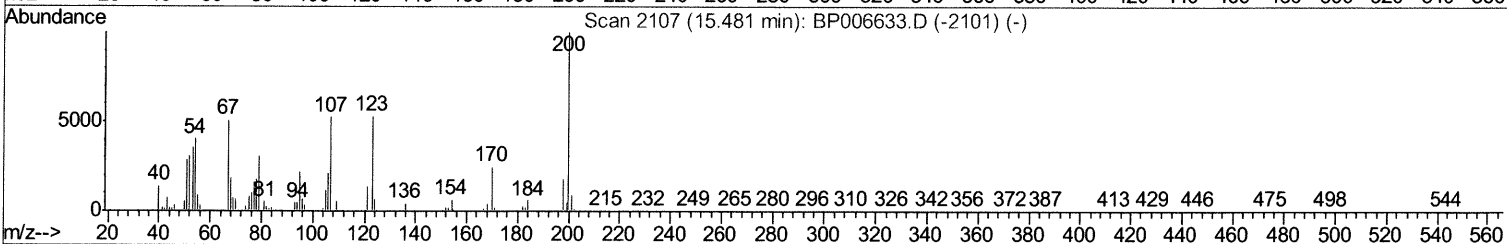
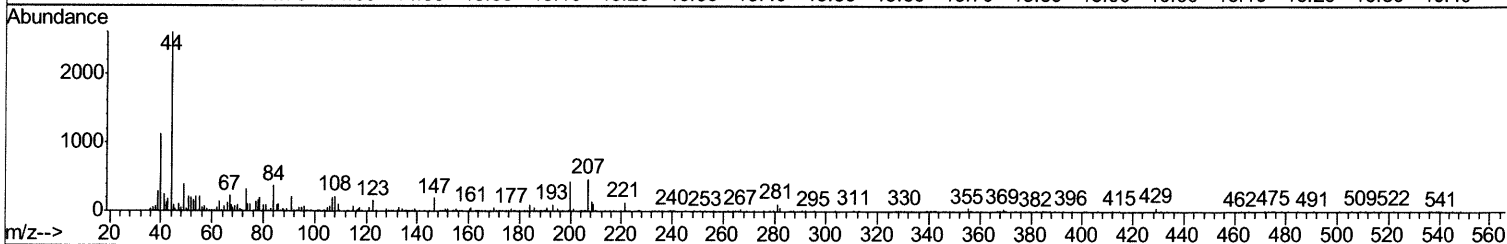
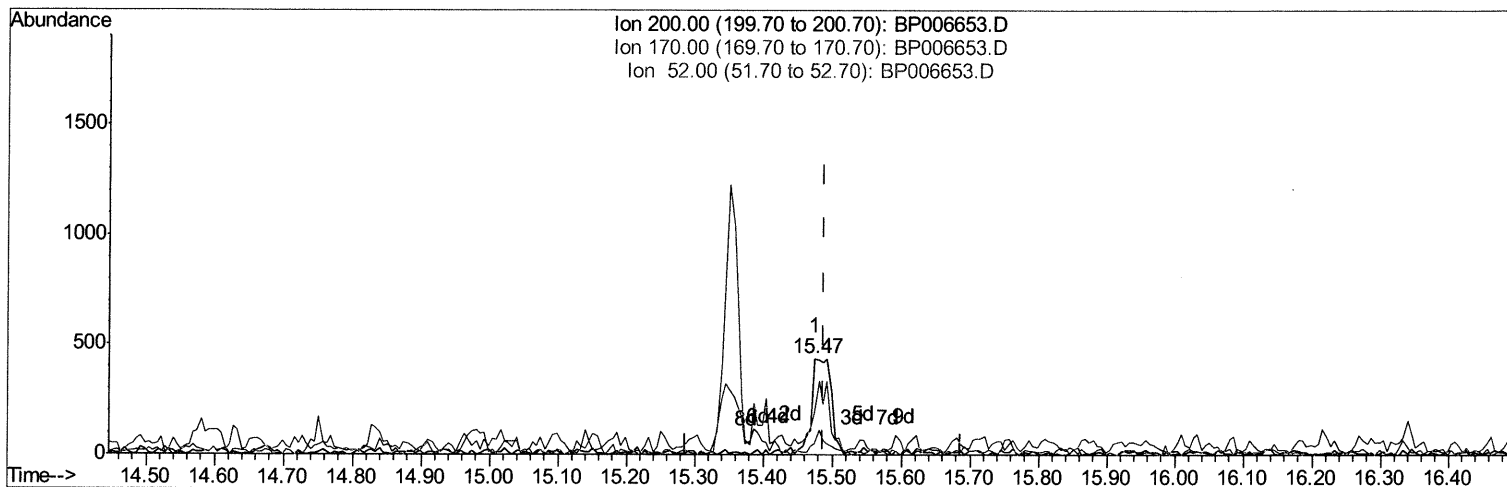
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TIC: BP006653.D

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.475min (-0.012) 0.19ng/ul m

response 851

Ion	Exp%	Act%
200.00	100	100
170.00	22.40	11.67#
52.00	47.50	43.71
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	145563	20.00	ng/ul	0.00
20) Naphthalene-d8	10.50	136	629084	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.36	164	359392	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.11	188	693843	20.00	ng/ul	0.00
79) Chrysene-d12	21.22	240	652188	20.00	ng/ul	0.00
88) Perylene-d12	23.53	264	674505	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	998m	0.23	ng/uL	0.00
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	0.00
7) Phenol-d5	6.90	99	6301	0.43	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.06	67	5294	0.59	ng/ul	0.00
11) 2-Chlorophenol-d4	7.25	132	5637	0.52	ng/ul	0.00
15) 4-Methylphenol-d8	8.43	113	3981	0.34	ng/ul	0.00
21) Nitrobenzene-d5	8.88	128	3046	0.57	ng/ul	0.00
24) 2-Nitrophenol-d4	9.59	143	2551	0.46	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.13	165	4659	0.49	ng/ul	0.00
31) 4-Chloroaniline-d4	10.65	131	4940	0.32	ng/ul	0.00
46) Dimethylphthalate-d6	13.78	166	15818	0.57	ng/ul	0.00
49) Acenaphthylene-d8	14.05	160	18433	0.55	ng/ul	0.00
54) 4-Nitrophenol-d4	14.59	143	650	0.11	ng/ul	0.01
60) Fluorene-d10	15.35	176	13630	0.61	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.47	200	851m	0.19	ng/ul	0.01
73) Anthracene-d10	17.21	188	21329	0.65	ng/ul	0.00
81) Pyrene-d10	19.46	212	23606	0.69	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.37	264	23186	0.64	ng/ul	0.00

Target Compounds Ovalue

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