

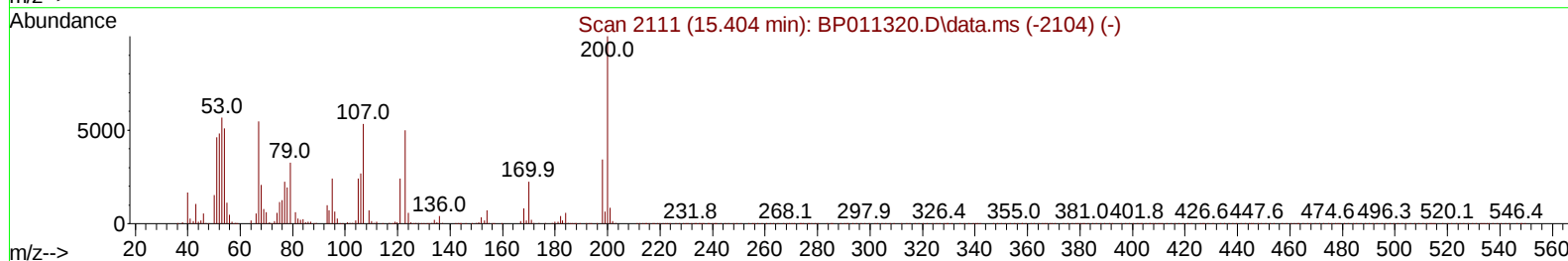
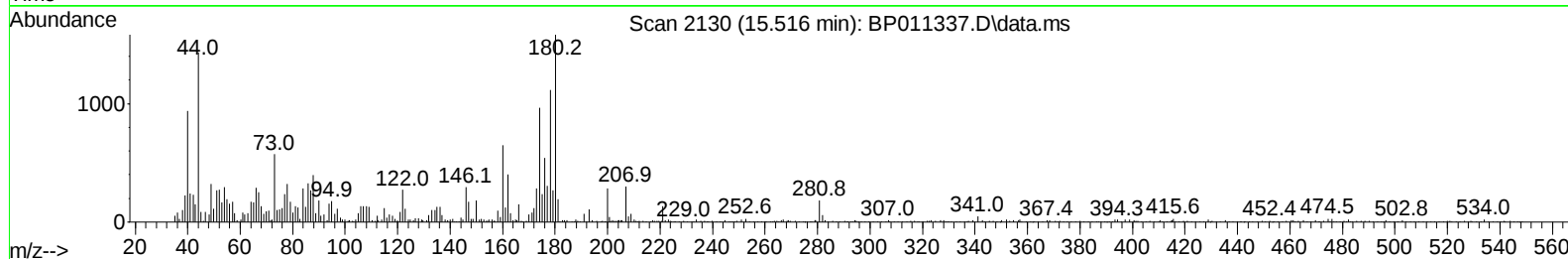
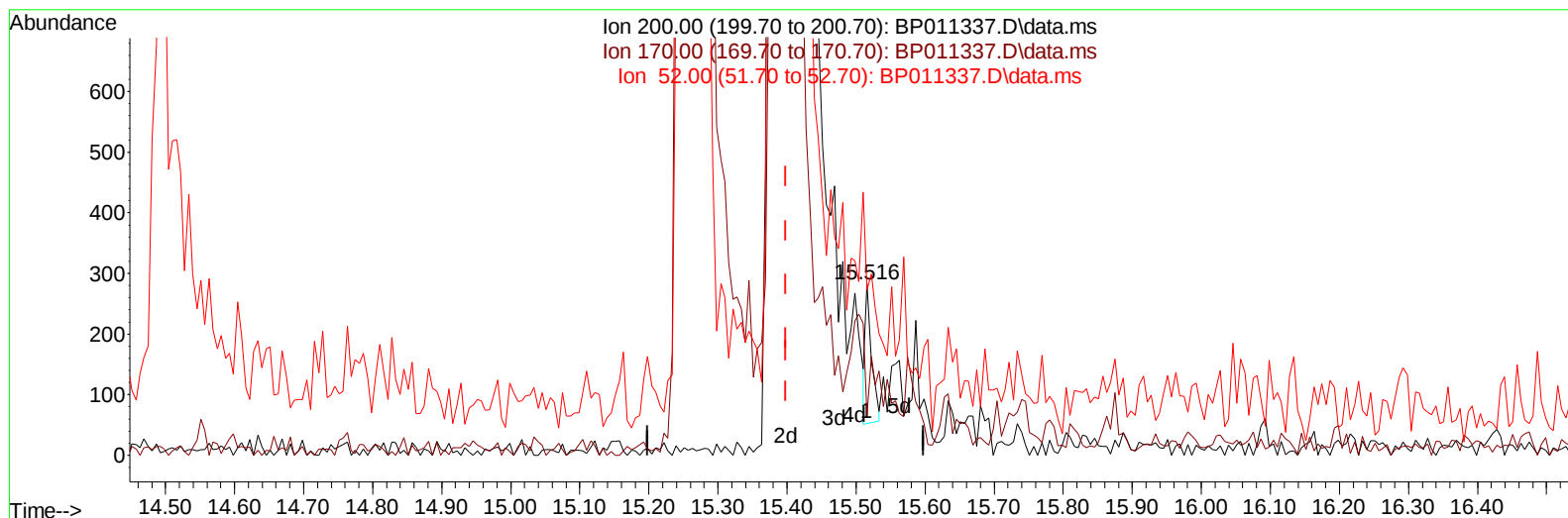
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011337.D
Acq On : 09 Aug 2022 19:34
Operator : CG/JU
Sample : N3900-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0C90

Manual IntegrationsAPPROVED

Quant Time: Aug 10 09:16:39 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/12/2022



TIC: BP011337.D\data.ms

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

15.516min (+ 0.118) 0.05 ng/u1

response 145

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.40	20.37
52.00	71.00	84.26
0.00	0.00	0.00

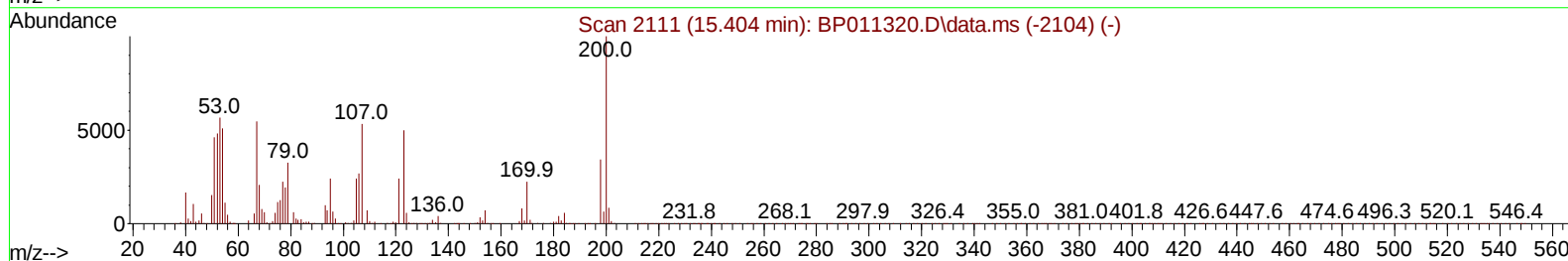
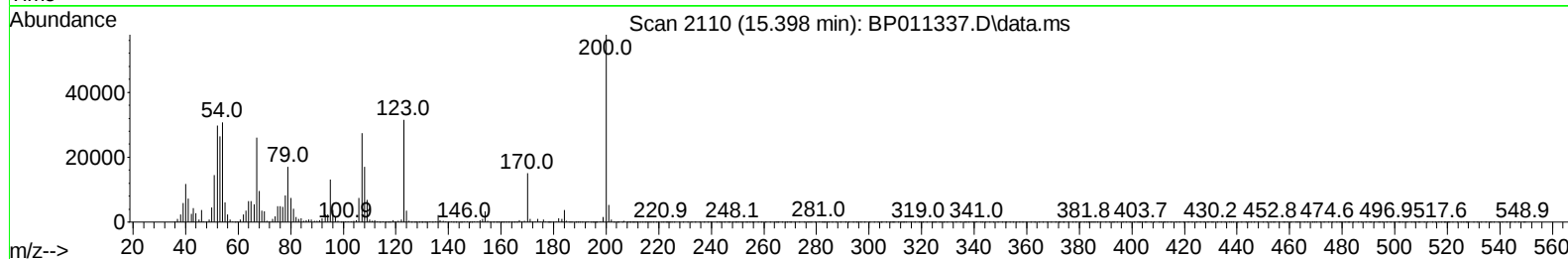
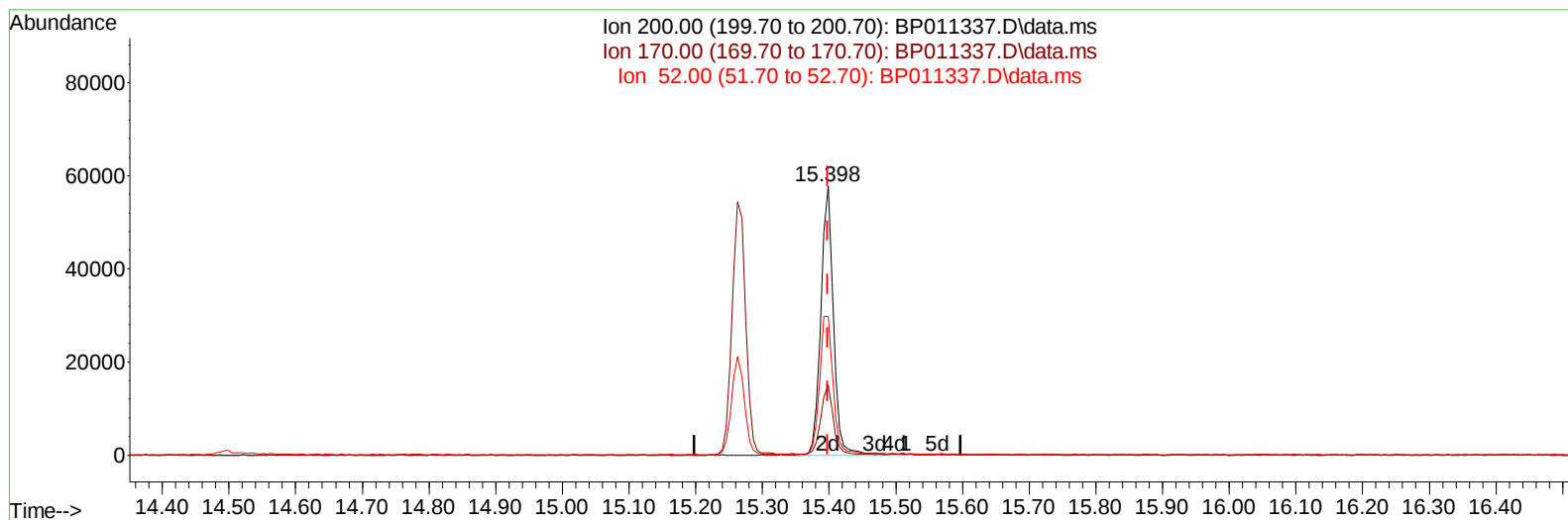
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.398min (-0.000) 26.59 ng/ul m

response 75314

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.40	26.10#
52.00	71.00	51.50#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	124360	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	530784	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	306237	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.034	188	609830	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	557120	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	574462	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	17001	4.191	ng/uL	0.00
4) Pyridine-d5	3.517	84	51209	5.004	ng/ul	0.02
7) Phenol-d5	6.775	99	53736	4.883	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	221172	28.343	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	174880	20.252	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	105854	12.250	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	118527	32.811	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	99719	30.151	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	164798	24.674	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	235861	20.199	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	716164	33.705	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	810180	31.675	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	9499	2.448	ng/ul	0.00
60) Fluorene-d10	15.263	176	593222	32.646	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	75314m	26.588	ng/ul	0.00
73) Anthracene-d10	17.133	188	1002657	36.579	ng/ul	0.00
81) Pyrene-d10	19.392	212	1137765	37.068	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	1063722	36.717	ng/ul	0.00

Target Compounds Qvalue

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

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