

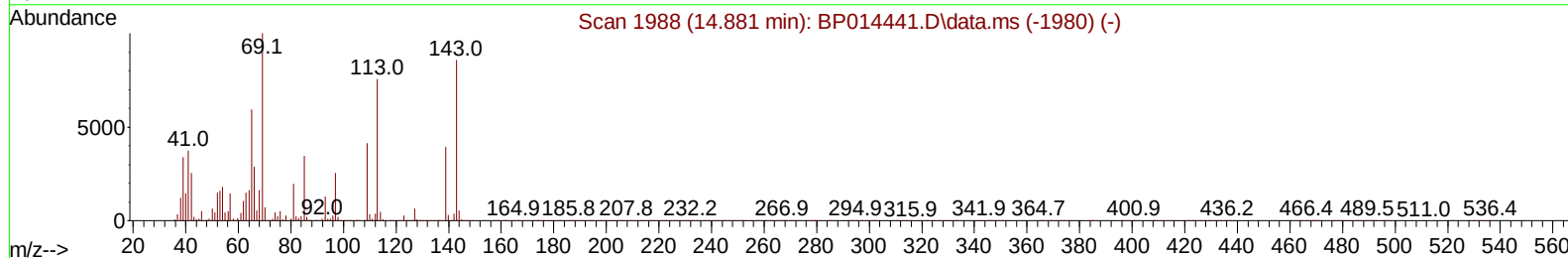
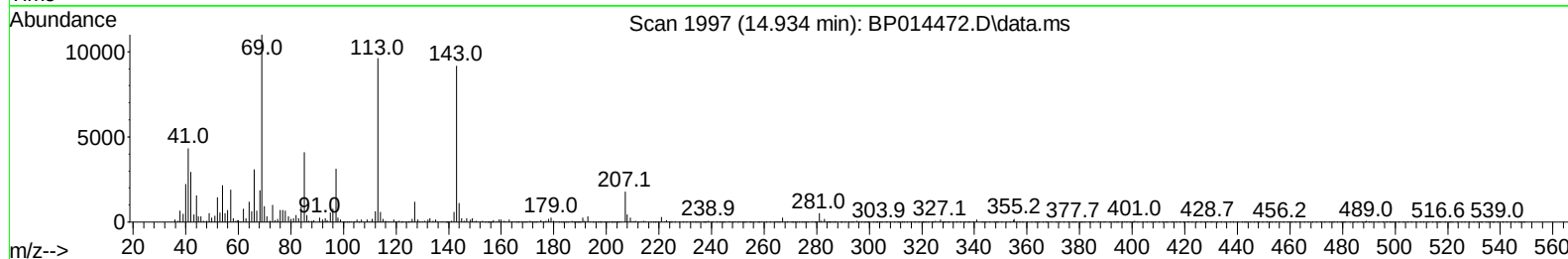
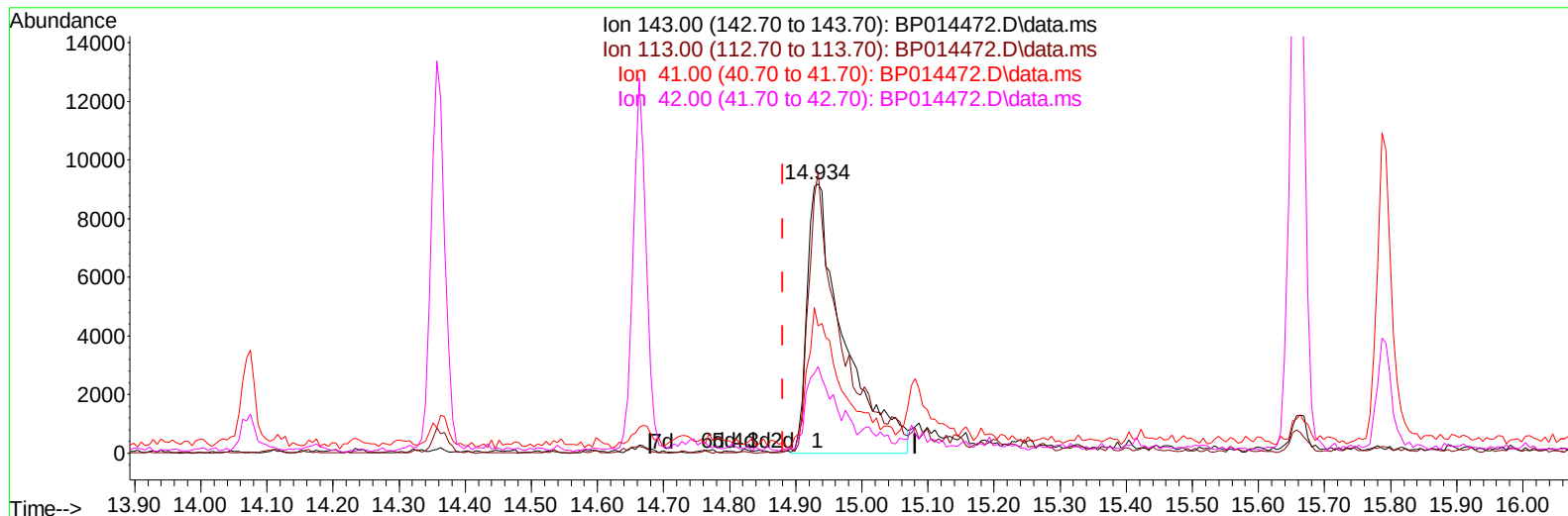
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
 Data File : BP014472.D
 Acq On : 01 Apr 2023 07:44
 Operator : CG/JU
 Sample : 02158-03
 Misc :
 ALS Vial : 82 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-10(20230329)

Manual IntegrationsAPPROVED

Quant Time: Apr 03 00:42:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 31 04:41:29 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/03/2023
 Supervised By :Jagrut Upadhyay 04/03/2023



TIC: BP014472.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.053) 5.65 ng/ul m

response 35397

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	104.62
41.00	44.20	47.31
42.00	31.50	32.17

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	155661	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	729656	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.663	164	511221	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1122666	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	857947	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	799972	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17225	4.291	ng/uL	0.00
4) Pyridine-d5	3.828	84	77311	7.516	ng/ul	0.01
7) Phenol-d5	7.187	99	175642	12.235	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	351891	38.204	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	300847	28.995	ng/ul	0.00
15) 4-Methylphenol-d8	8.740	113	292551	25.264	ng/ul	0.01
21) Nitrobenzene-d5	9.187	128	176131	29.901	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	201790	29.917	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	361792	29.453	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	378803	23.411	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	1321727	31.887	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	1398394	30.404	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143	35397m	5.653	ng/ul	0.05
60) Fluorene-d10	15.663	176	1106596	32.013	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	184131	22.290	ng/ul	0.00
73) Anthracene-d10	17.528	188	1892389	34.830	ng/ul	0.00
81) Pyrene-d10	19.757	212	2128502	36.935	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	1534362	36.101	ng/ul	0.00

Target Compounds Qvalue

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

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