

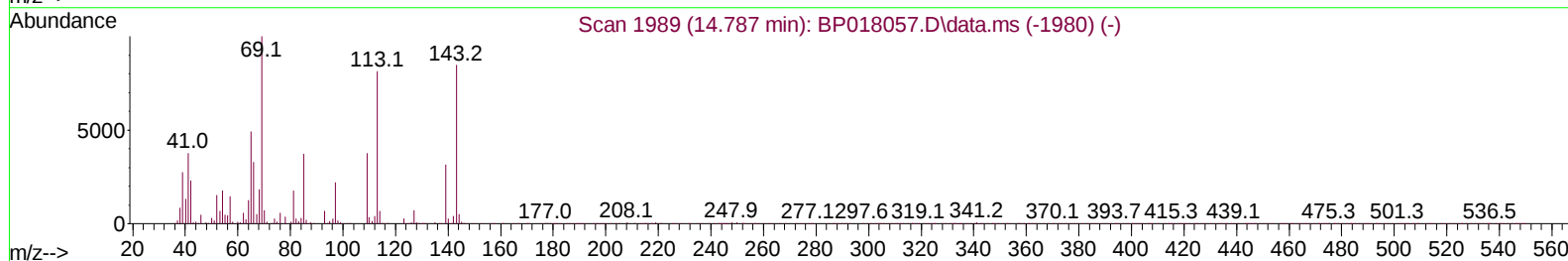
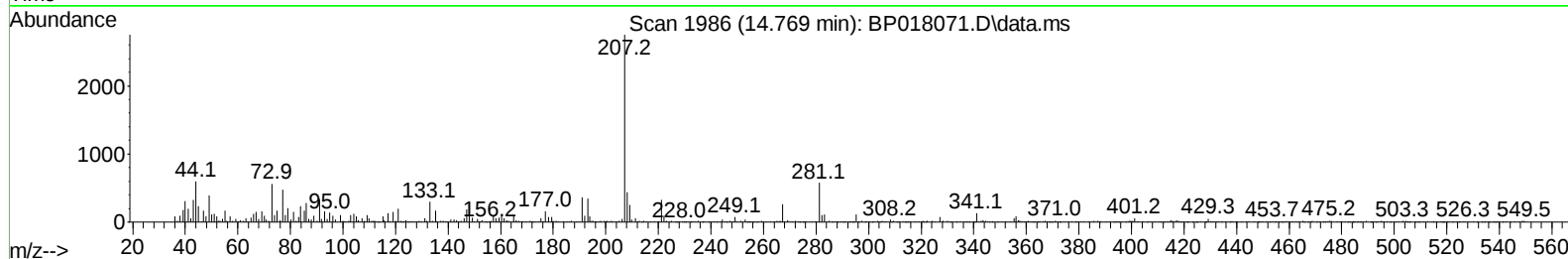
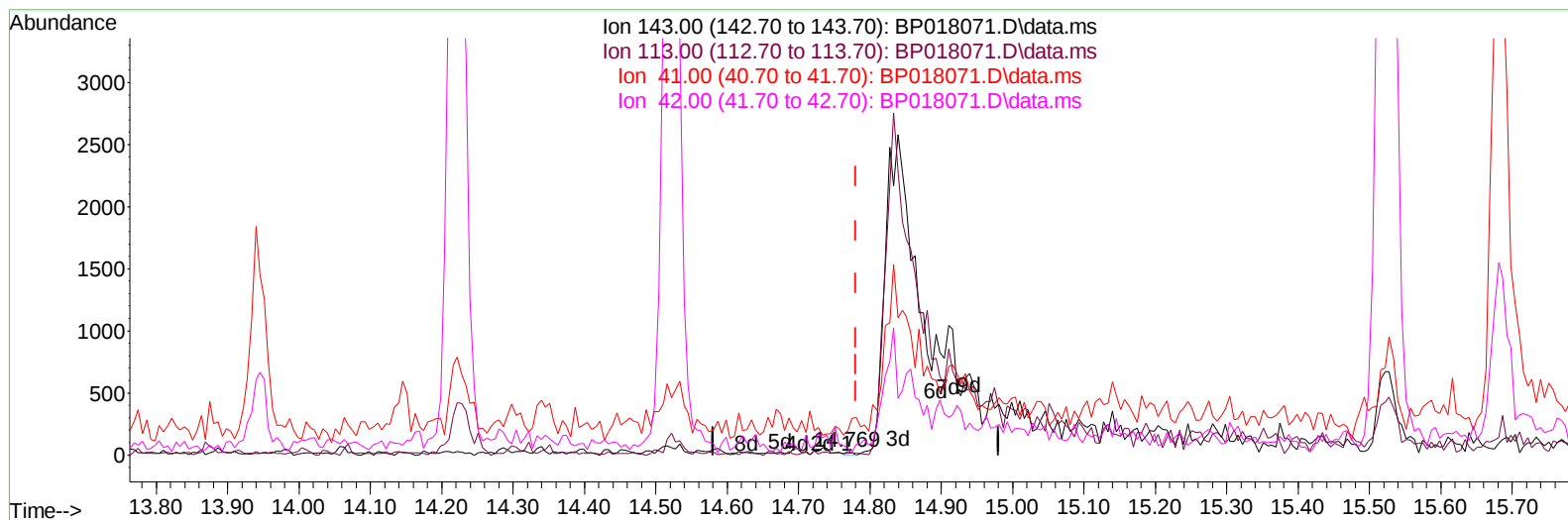
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018071.D
 Acq On : 11 Nov 2023 18:02
 Operator : MA/JU
 Sample : 05016-07
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A49A9

Manual IntegrationsAPPROVED

Quant Time: Nov 11 21:59:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/12/2023
 Supervised By :mohammad ahmed 11/15/2023



TIC: BP018071.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.012) 0.00 ng/ul

response 17

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	44.44#
41.00	47.70	733.33#
42.00	30.10	196.30#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/12/2023
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Quant Time: Nov 11 22:00:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	94749	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	404557	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	248025	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	473774	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	366991	20.000	ng/ul	0.00
88) Perylene-d12	23.933	264	432637	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	8567	3.190	ng/uL	0.00
4) Pyridine-d5	3.669	84	66234	9.458	ng/ul	0.00
7) Phenol-d5	7.022	99	51881	5.672	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	153038	28.469	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	149516	20.641	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	96777	12.911	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	98235	26.952	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	106580	25.811	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	170199	23.126	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	199143	19.731	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	635113	27.753	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	666666	26.746	ng/ul	0.00
54) 4-Nitrophenol-d4	14.839	143	9539m	2.789	ng/ul	0.06
60) Fluorene-d10	15.528	176	502966	26.621	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	75723	22.109	ng/ul	0.00
73) Anthracene-d10	17.404	188	714259	27.875	ng/ul	0.00
81) Pyrene-d10	19.657	212	833721	34.011	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.768	264	802937	31.918	ng/ul	0.00

Target Compounds Qvalue

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

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