

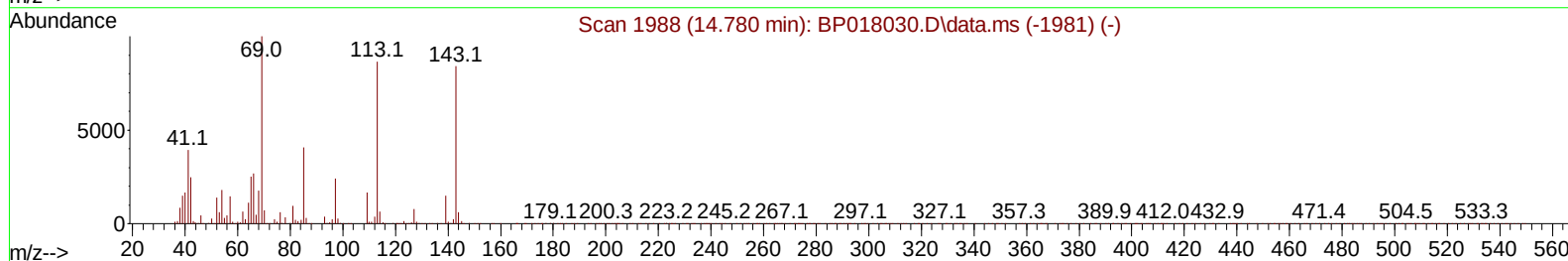
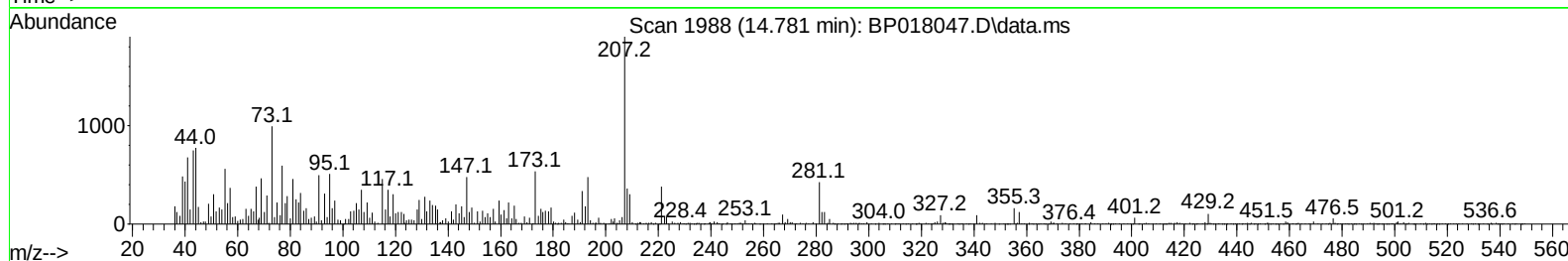
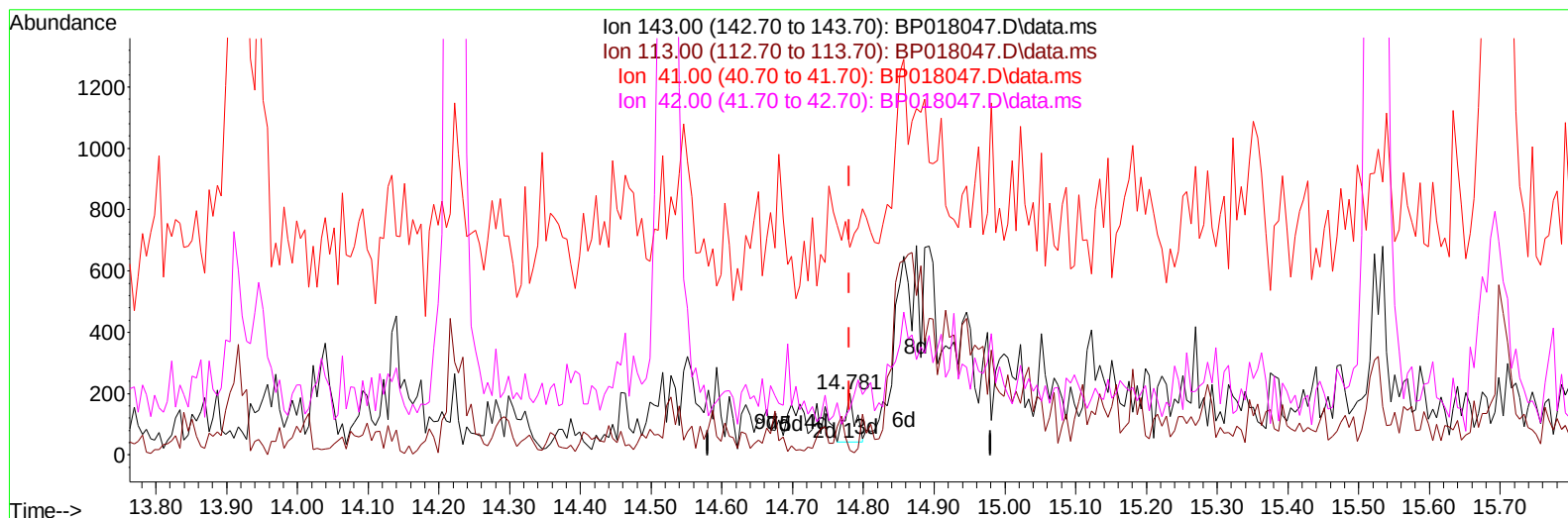
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018047.D
 Acq On : 11 Nov 2023 03:56
 Operator : MA/JU
 Sample : 05044-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHEE3

Manual IntegrationsAPPROVED

Quant Time: Nov 11 04:28:39 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: BP018047.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.781min (+ 0.000) 0.08 ng/ul

response 158

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	7.96#
41.00	47.70	336.32#
42.00	30.10	74.63#

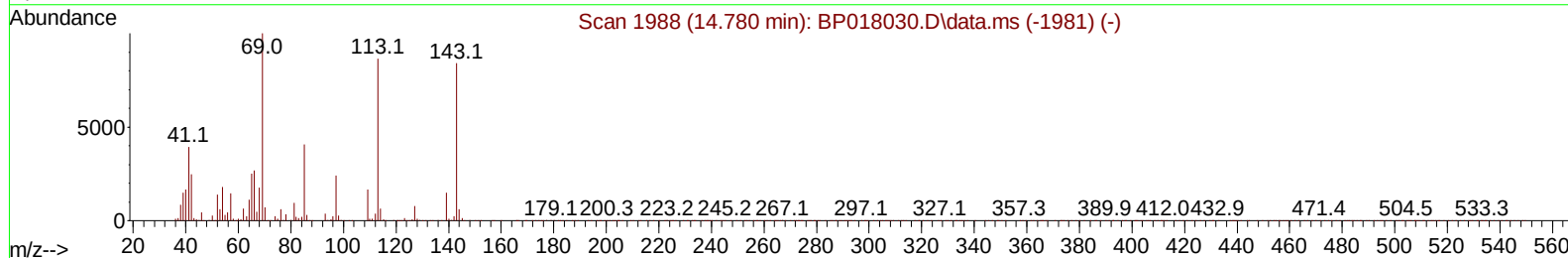
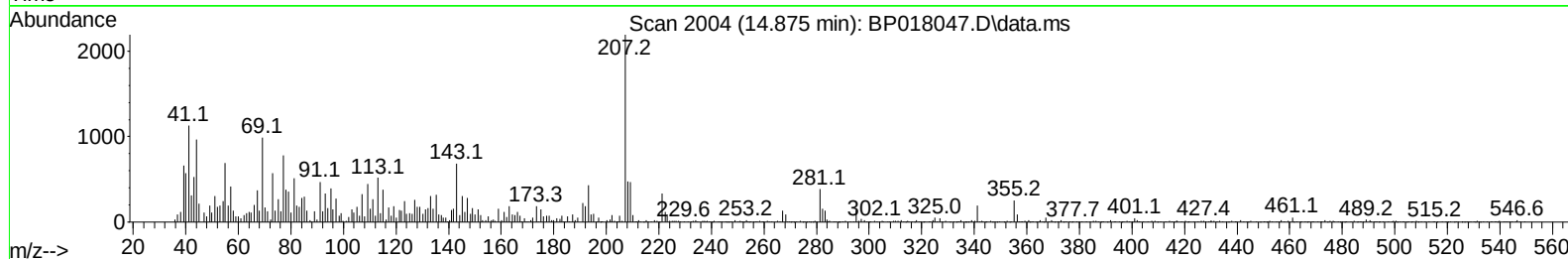
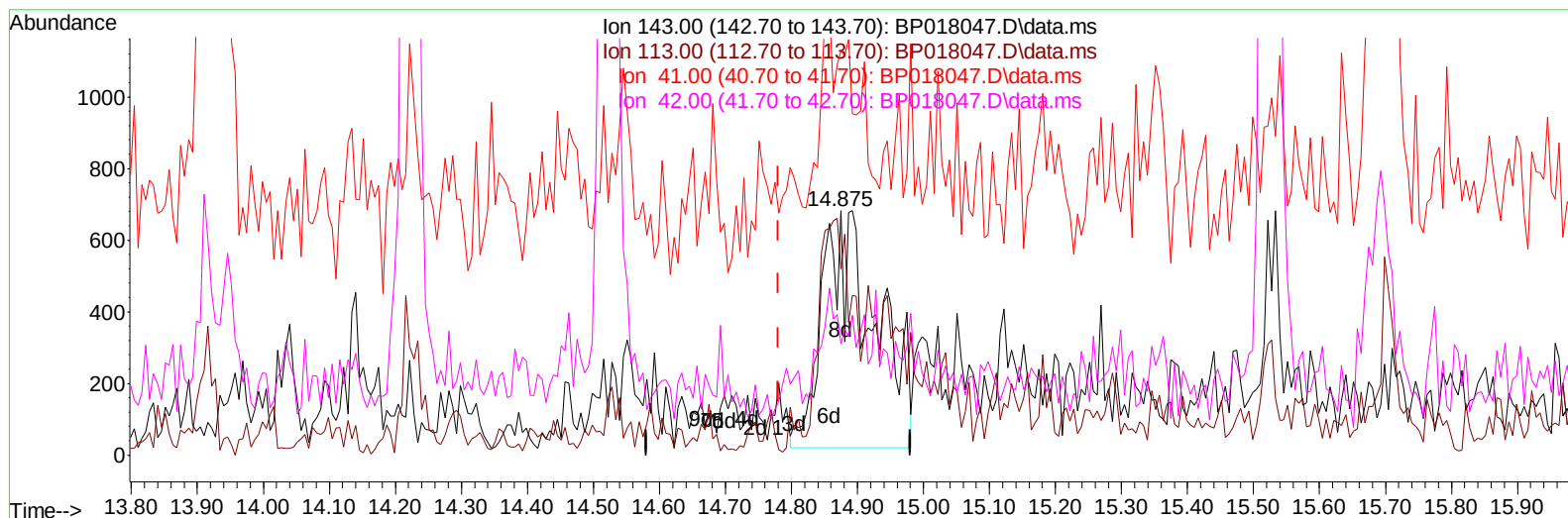
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Manual IntegrationsAPPROVED

Quant Time: Nov 11 04:29:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
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TIC: BP018047.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.094) 1.80 ng/ul m

response 3680

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	76.13#
41.00	47.70	165.30#
42.00	30.10	45.53#

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 Operator : MA/JU
 Sample : 05044-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHEE3

Manual IntegrationsAPPROVED

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

Quant Time: Nov 11 04:29:36 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	65796	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	255347	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	147999	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	306036	20.000	ng/ul	0.00
79) Chrysene-d12	21.427	240	275992	20.000	ng/ul	0.00
88) Perylene-d12	23.933	264	329519	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	5153	2.763	ng/uL	0.00
4) Pyridine-d5	3.670	84	45405	9.336	ng/ul	0.00
7) Phenol-d5	7.034	99	17549	2.763	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.199	67	92656	24.821	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	59319	11.792	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113	41472	7.967	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	56825	24.701	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	40580	15.570	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	63732	13.720	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	57432	9.016	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	363743	26.637	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	394715	26.539	ng/ul	0.00
54) 4-Nitrophenol-d4	14.875	143	3680m	1.803	ng/ul	0.09
60) Fluorene-d10	15.528	176	305801	27.124	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	25769	11.647	ng/ul	0.00
73) Anthracene-d10	17.404	188	493226	29.799	ng/ul	0.00
81) Pyrene-d10	19.657	212	568816	30.855	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.769	264	577066	30.117	ng/ul	0.00

Target Compounds Qvalue

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

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