

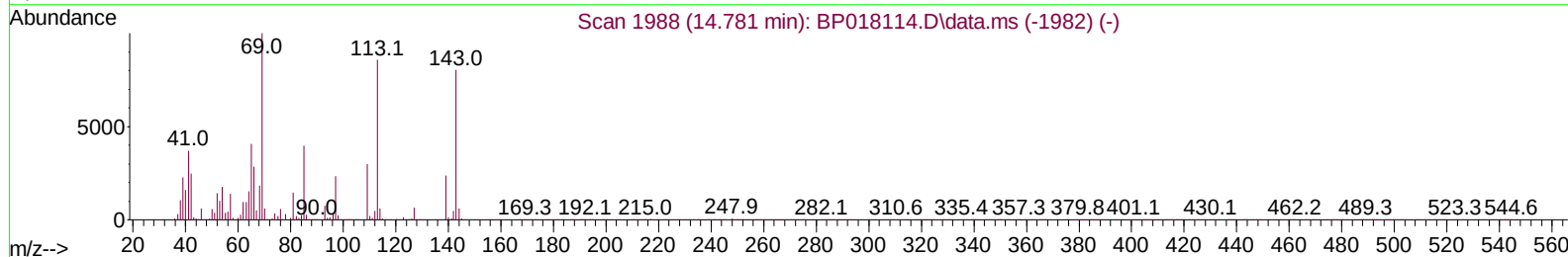
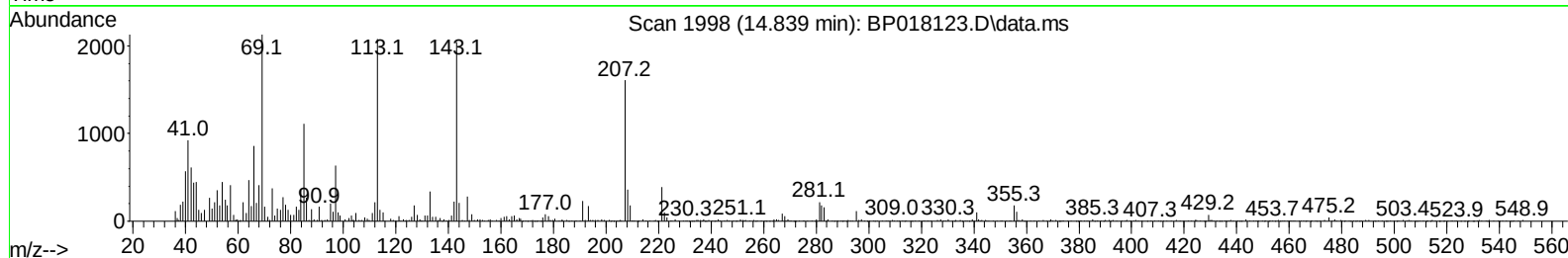
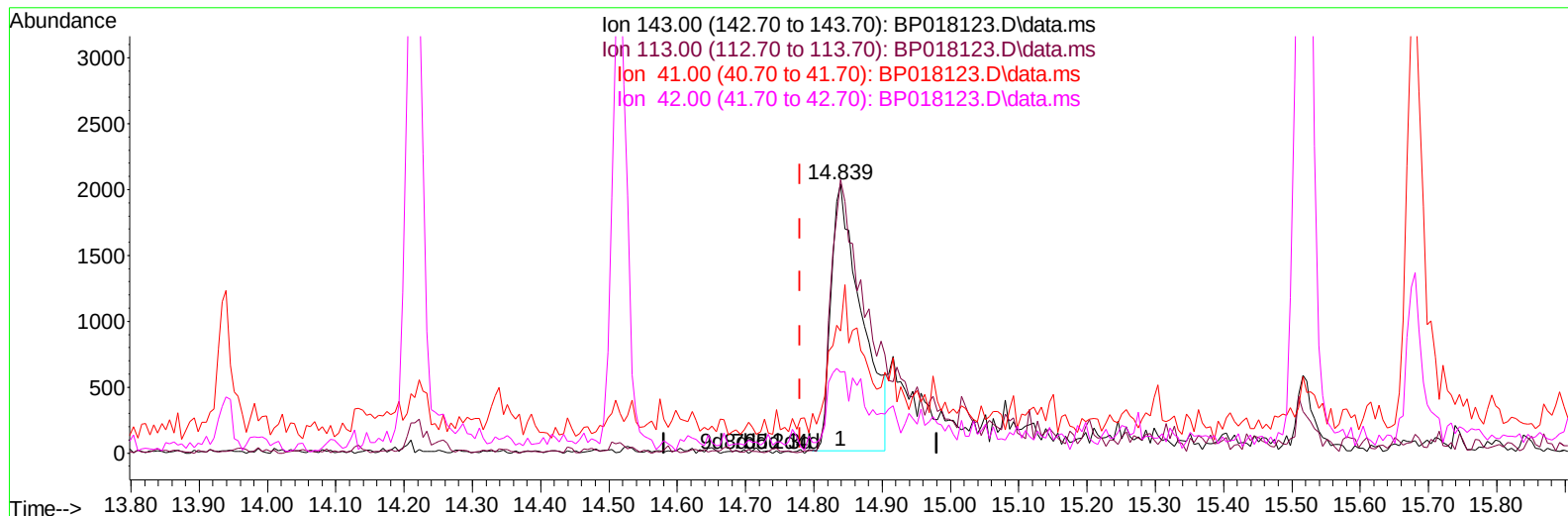
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH362

Manual IntegrationsAPPROVED

Quant Time: Nov 13 03:40:37 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 11/14/2023
Supervised By :mohammad ahmed 11/15/2023



TIC: BP018123.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.839min (+ 0.059) 4.22 ng/u1

response 6256

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	101.86
41.00	47.70	45.42
42.00	30.10	29.99

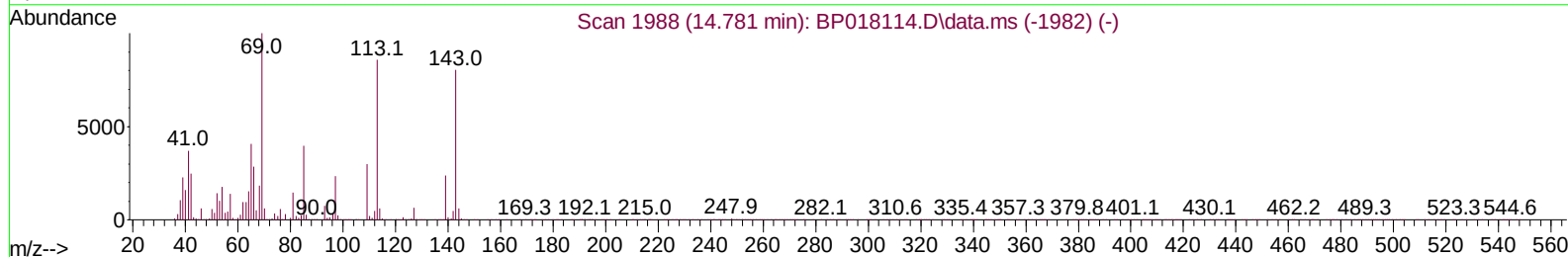
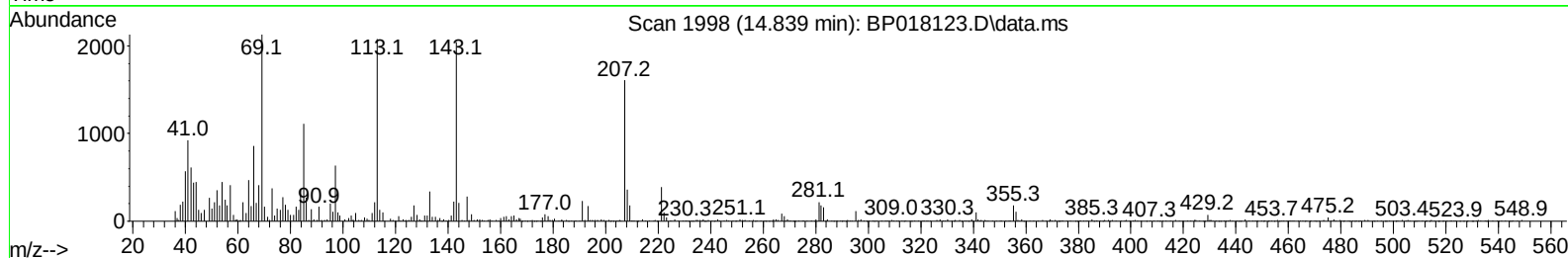
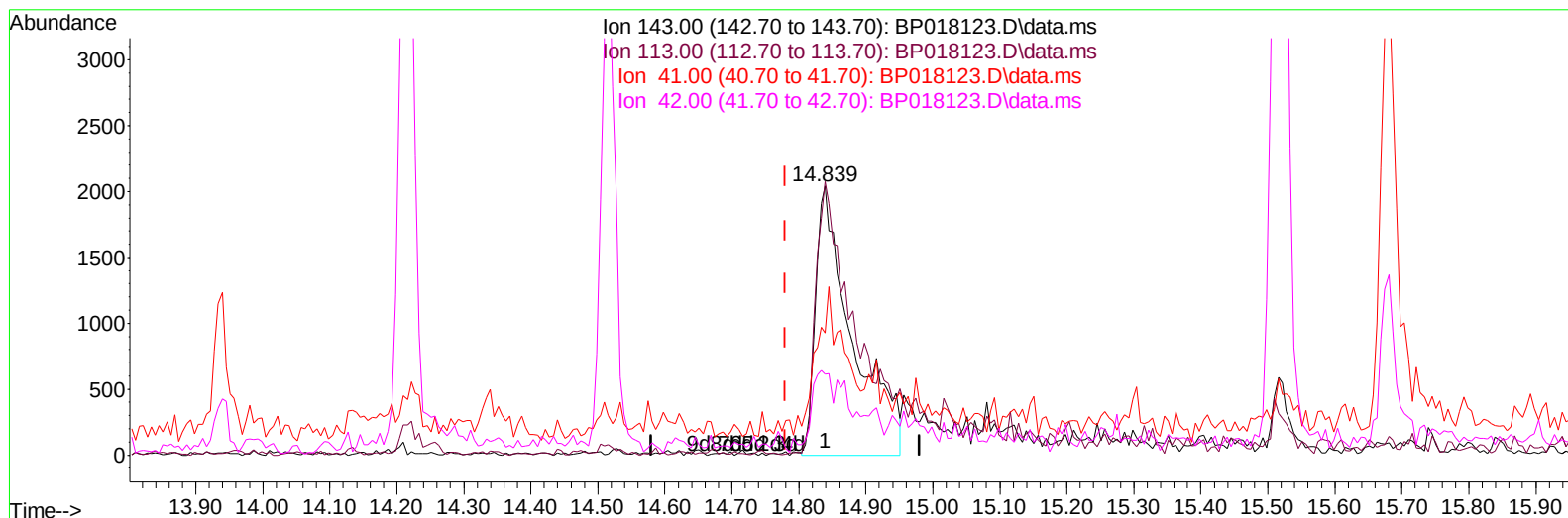
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TIC: BP018123.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.839min (+ 0.059) 5.26 ng/ul m

response 7800

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	101.86
41.00	47.70	45.42
42.00	30.10	29.99

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Manual IntegrationsAPPROVED

Quant Time: Nov 13 03:41:48 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 11/14/2023
 Supervised By :mohammad ahmed 11/15/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	47097	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.663	136	185611	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	107602	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.298	188	242445	20.000	ng/ul	-0.01
79) Chrysene-d12	21.427	240	266115	20.000	ng/ul	0.00
88) Perylene-d12	23.921	264	340876	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	6464	4.842	ng/uL	0.00
4) Pyridine-d5	3.669	84	37035	10.639	ng/ul	0.00
7) Phenol-d5	7.022	99	33615	7.393	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	95107	35.593	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	96140	26.701	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	61659	16.549	ng/ul	0.00
21) Nitrobenzene-d5	9.051	128	57209	34.211	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	63720	33.634	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.292	165	104903	31.068	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	134914	29.136	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	383465	38.624	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	394323	36.466	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.839	143	7800m	5.258	ng/ul	0.06
60) Fluorene-d10	15.522	176	320803	39.137	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.680	200	53107	30.300	ng/ul	0.00
73) Anthracene-d10	17.398	188	555364	42.354	ng/ul	-0.01
81) Pyrene-d10	19.651	212	725881	40.836	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.756	264	858500	43.313	ng/ul	-0.02
Target Compounds						
78) Di-n-butylphthalate	18.233	149	17447	1.042	ng/ul#	95

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

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