

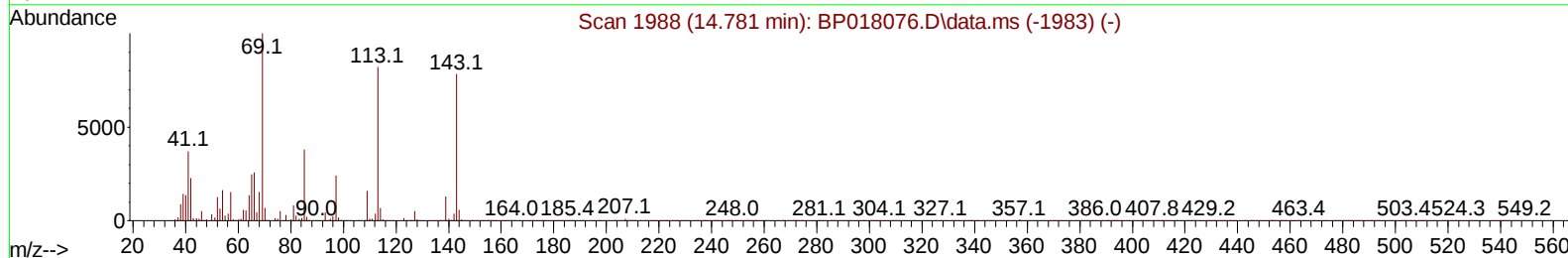
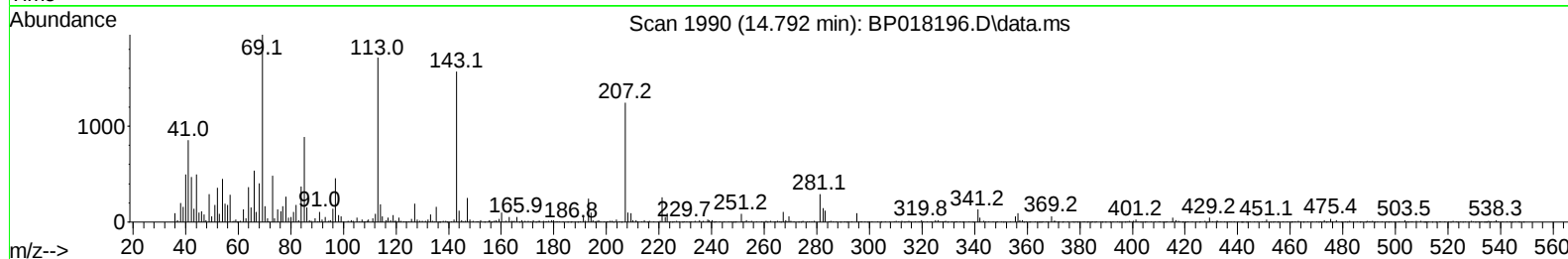
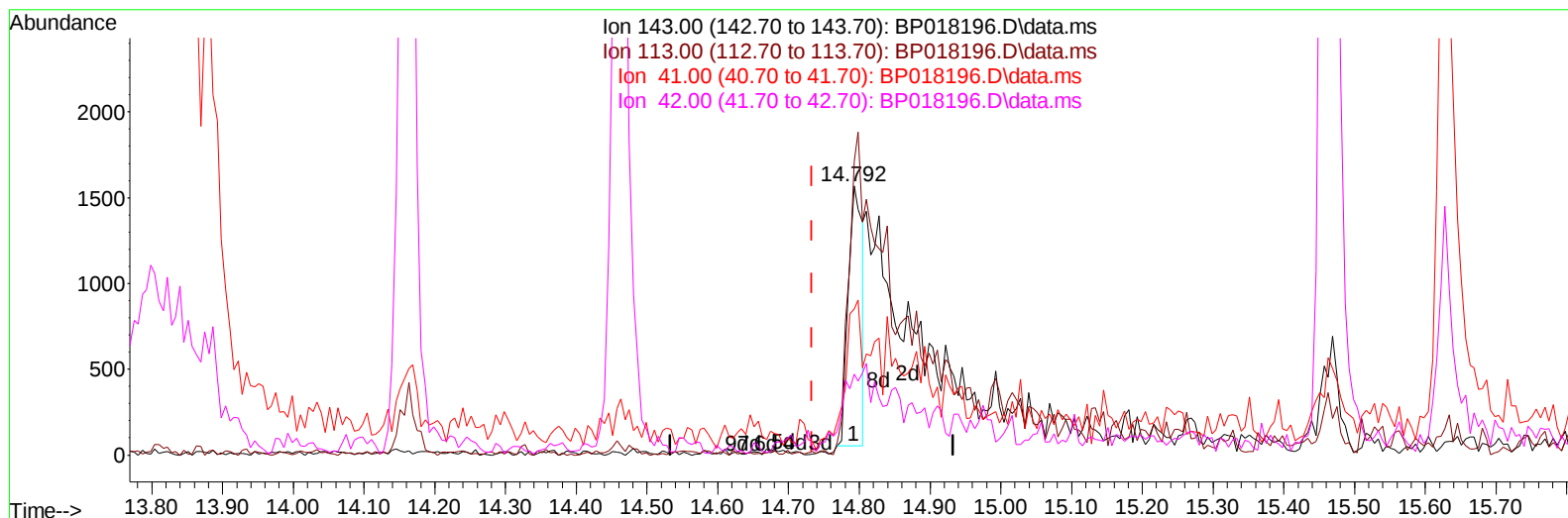
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018196.D
 Acq On : 16 Nov 2023 04:25
 Operator : MA/JU
 Sample : 05338-20
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDM5

Manual IntegrationsAPPROVED

Quant Time: Nov 16 05:00:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
 Supervised By :mohammad ahmed 11/17/2023



TIC: BP018196.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.792min (+ 0.059) 0.96 ng/u1

response 2071

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	109.25
41.00	47.70	54.40
42.00	30.10	30.17

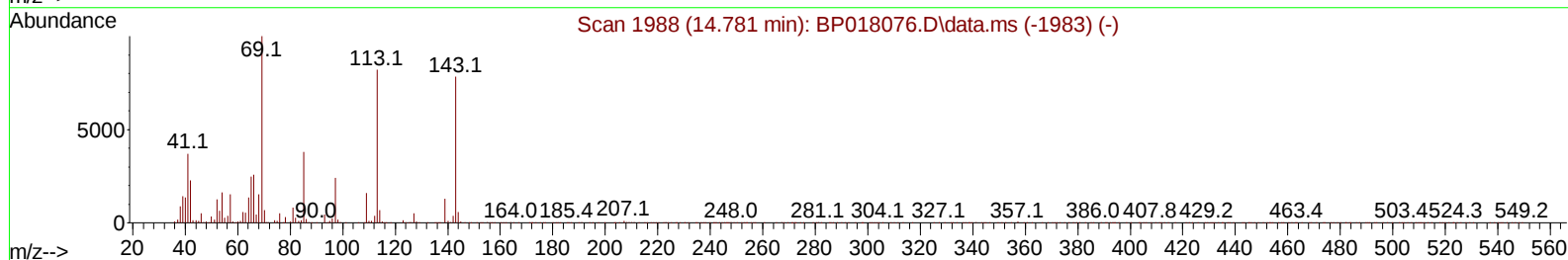
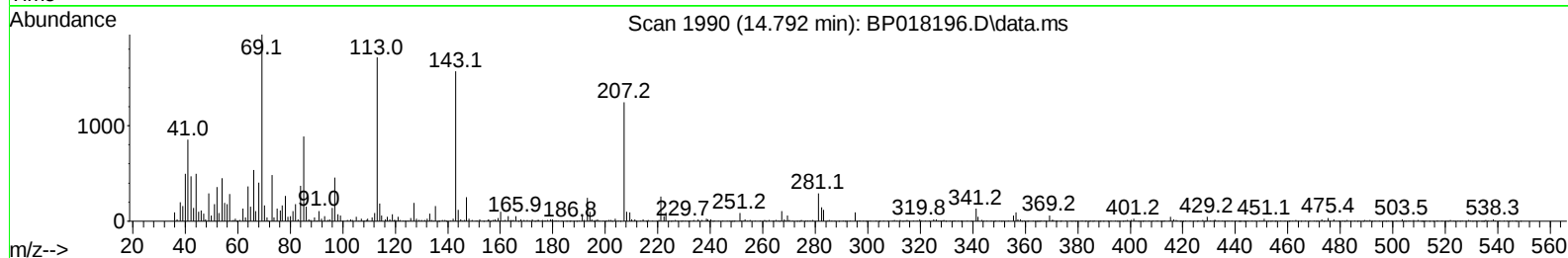
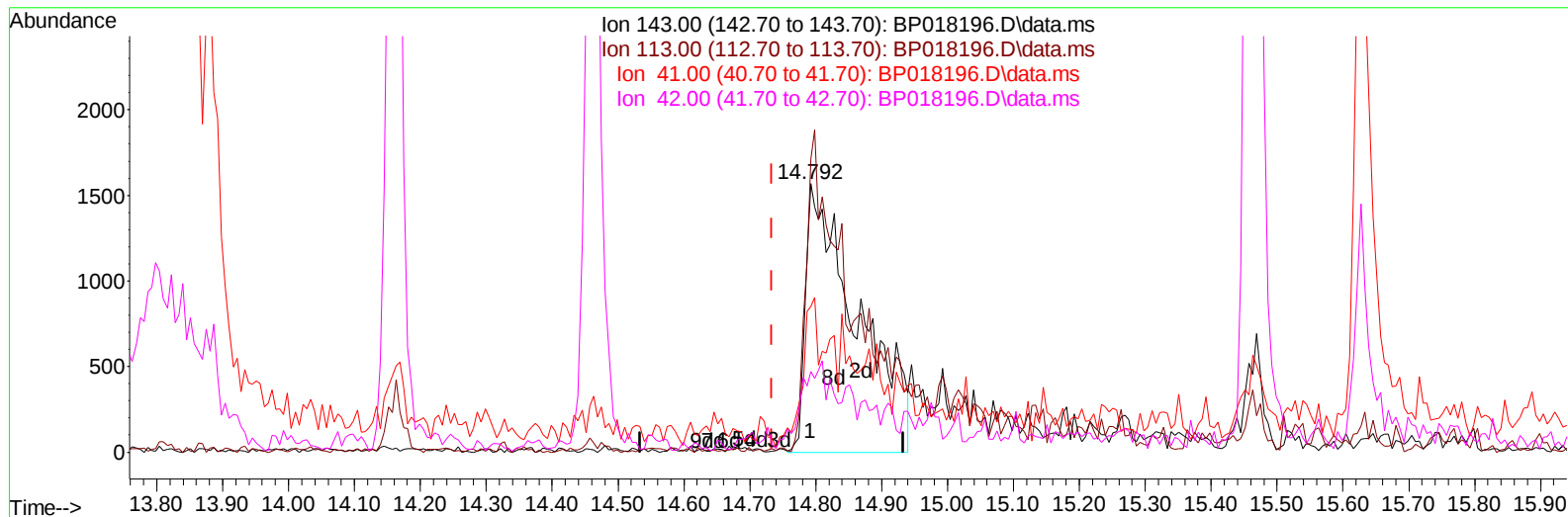
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TIC: BP018196.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.792min (+ 0.059) 3.93 ng/ul m

response 8511

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	109.25
41.00	47.70	54.40
42.00	30.10	30.17

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	71965	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	268869	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	157211	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	340051	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	349283	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	404336	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	8068	3.955	ng/uL	0.00
4) Pyridine-d5	3.605	84	34707	6.525	ng/ul	0.01
7) Phenol-d5	6.963	99	41567	5.983	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	117020	28.660	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	121144	22.019	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	75085	13.189	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	68246	28.173	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	72998	26.600	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	113932	23.293	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	93706	13.970	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	440179	30.346	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	453550	28.707	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	8511m	3.927	ng/ul	0.06
60) Fluorene-d10	15.463	176	356636	29.779	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	55761	22.683	ng/ul	0.00
73) Anthracene-d10	17.345	188	595743	32.393	ng/ul	0.00
81) Pyrene-d10	19.598	212	747605	32.044	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	781425	33.237	ng/ul	0.00

Target Compounds Qvalue

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

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