

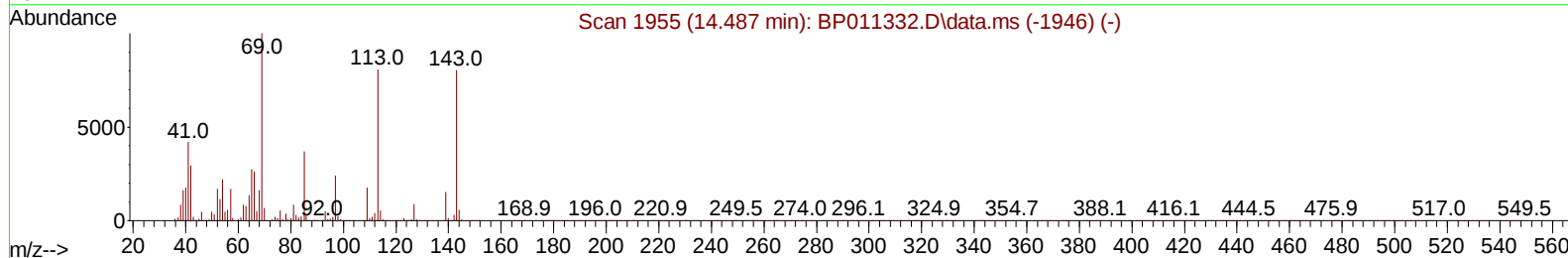
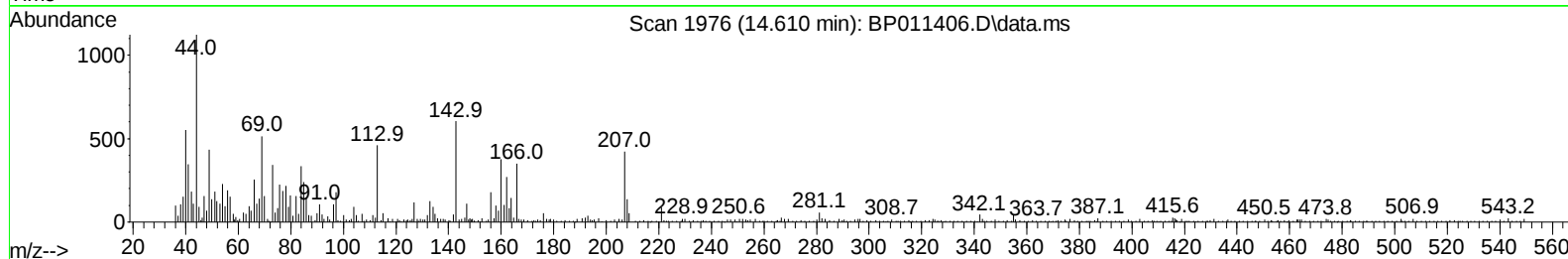
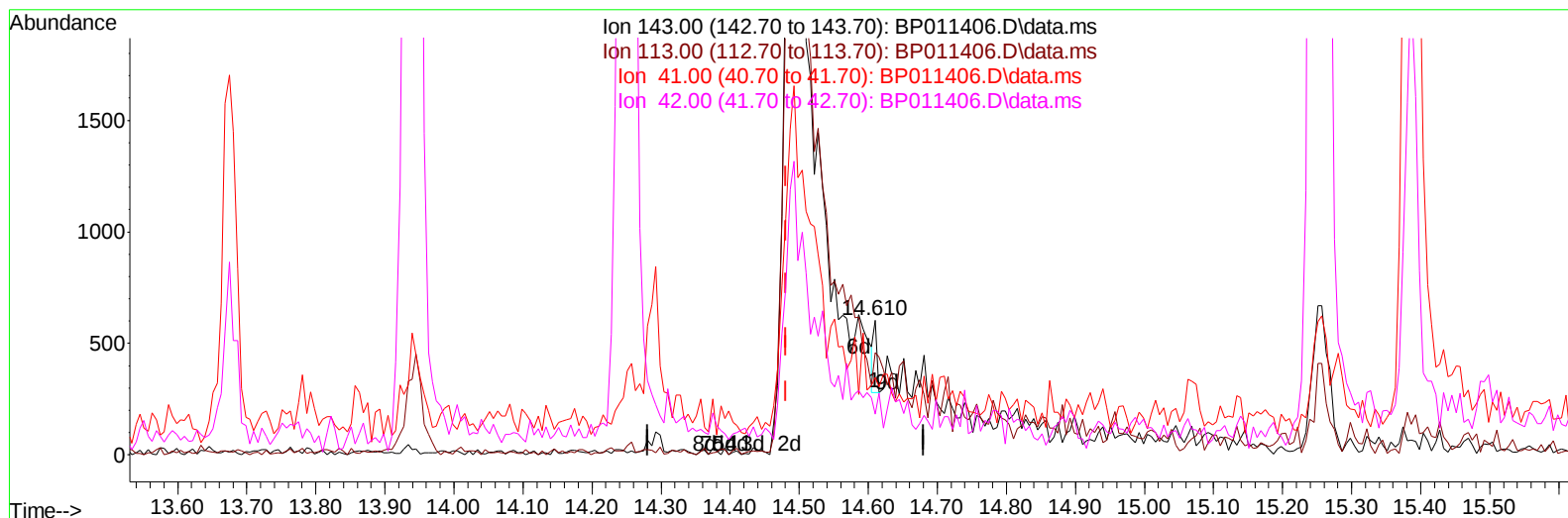
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011406.D
 Acq On : 12 Aug 2022 02:33
 Operator : CG/JU
 Sample : N4128-05
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBW6

Manual IntegrationsAPPROVED

Quant Time: Aug 12 03:20:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 11 23:05:53 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/12/2022
 Supervised By :Sohil Jodhani 08/13/2022



TIC: BP011406.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.129) 0.05 ng/u1

response 114

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	75.87
41.00	54.40	57.52
42.00	35.50	30.41

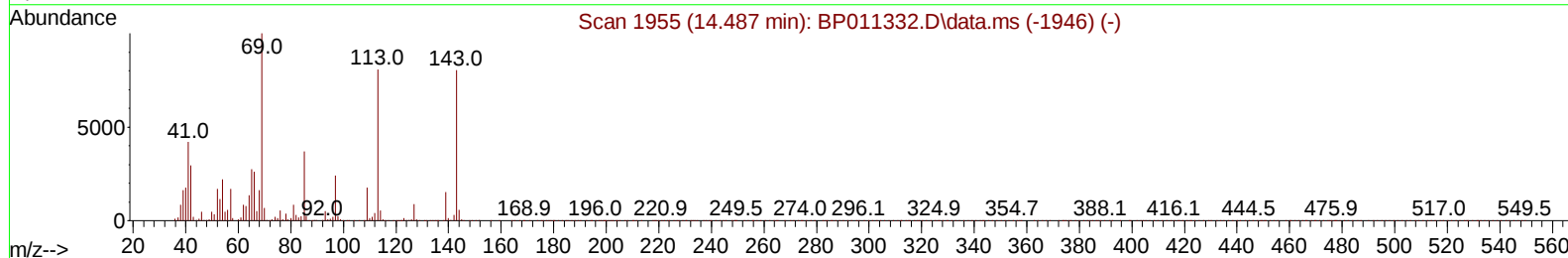
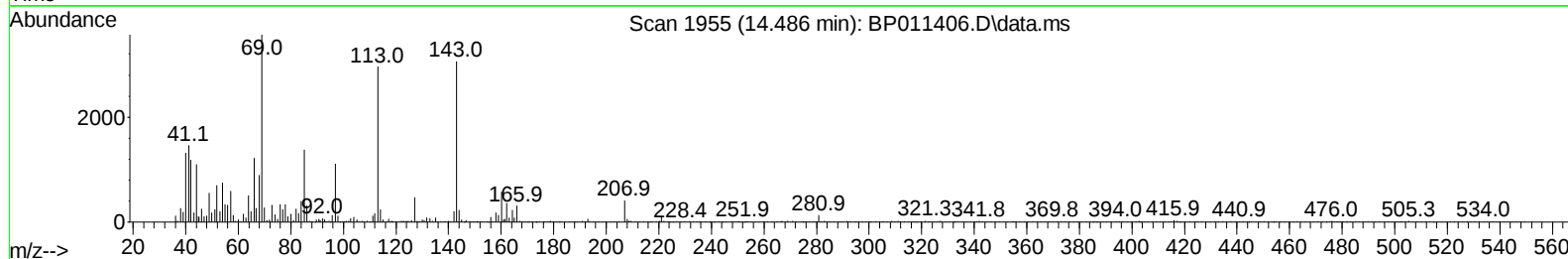
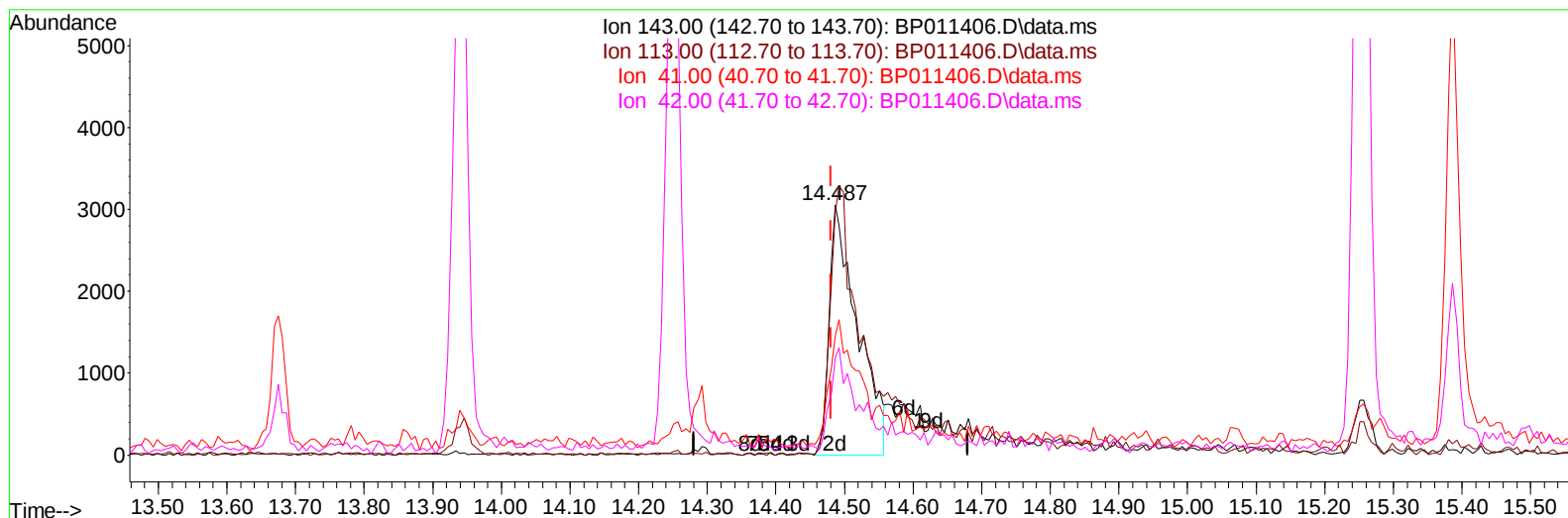
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TIC: BP011406.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.486min (+ 0.006) 3.57 ng/ul m

response 8803

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	96.82#
41.00	54.40	47.89
42.00	35.50	38.89

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	82120	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	345471	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	194401	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	410164	20.000	ng/ul	0.00
79) Chrysene-d12	21.157	240	400003	20.000	ng/ul	# 0.00
88) Perylene-d12	23.462	264	420600	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	10511	3.924	ng/uL	0.00
4) Pyridine-d5	3.499	84	50014	7.402	ng/ul	0.01
7) Phenol-d5	6.763	99	48335	6.651	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	146557	28.441	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	138837	24.348	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	86957	15.239	ng/ul	0.00
21) Nitrobenzene-d5	8.740	128	80143	34.086	ng/ul	0.00
24) 2-Nitrophenol-d4	9.457	143	76542	35.558	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	120667	27.758	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	170273	22.405	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	482143	35.745	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	537727	33.118	ng/ul	0.00
54) 4-Nitrophenol-d4	14.486	143	8803m	3.573	ng/ul	0.00
60) Fluorene-d10	15.257	176	394021	34.158	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	55960	29.373	ng/ul	0.00
73) Anthracene-d10	17.122	188	679309	36.847	ng/ul	0.00
81) Pyrene-d10	19.386	212	787857	35.751	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	777367	36.649	ng/ul	0.00

Target Compounds Qvalue

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

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