

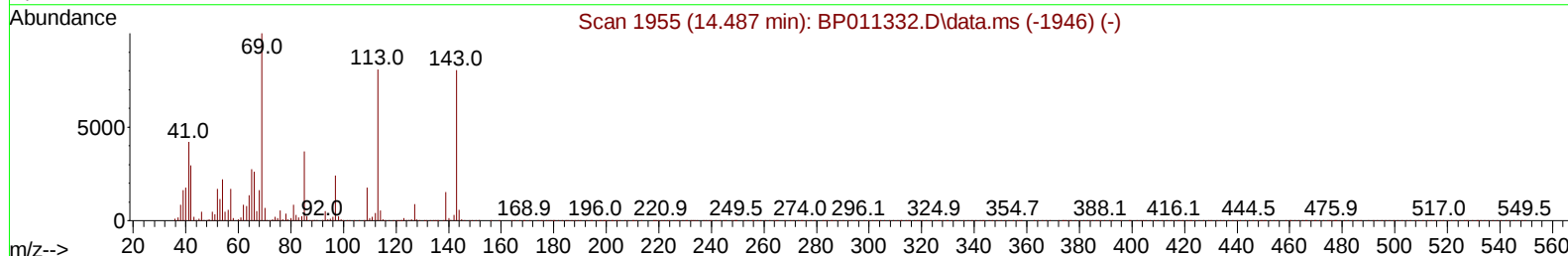
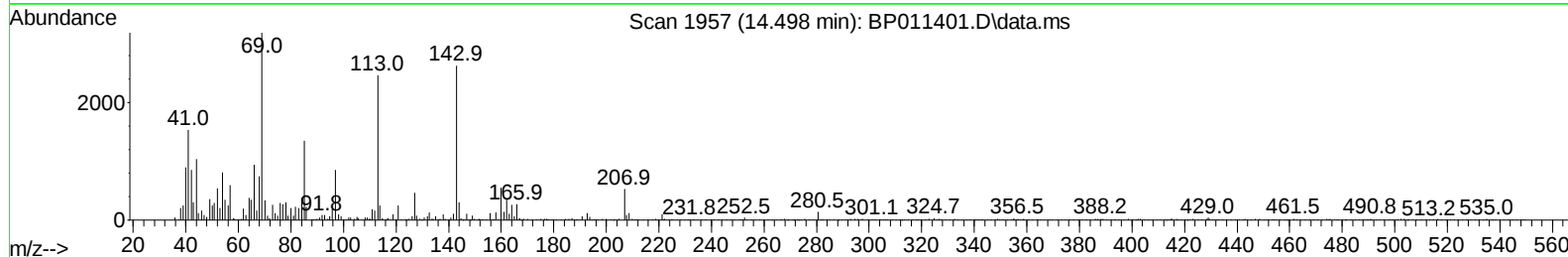
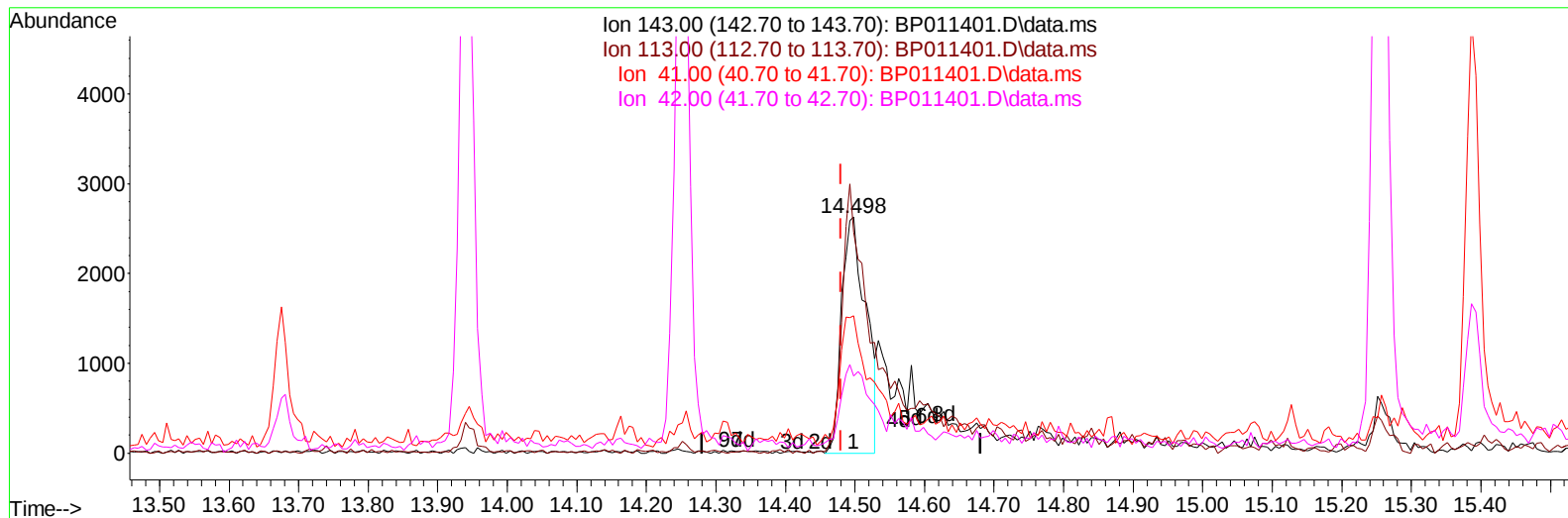
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011401.D
 Acq On : 11 Aug 2022 23:41
 Operator : CG/JU
 Sample : N4128-10
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBT7

Manual IntegrationsAPPROVED

Quant Time: Aug 12 01:10:55 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: BP011401.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 2.47 ng/ul

response 6366

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	93.50#
41.00	54.40	58.19
42.00	35.50	32.57

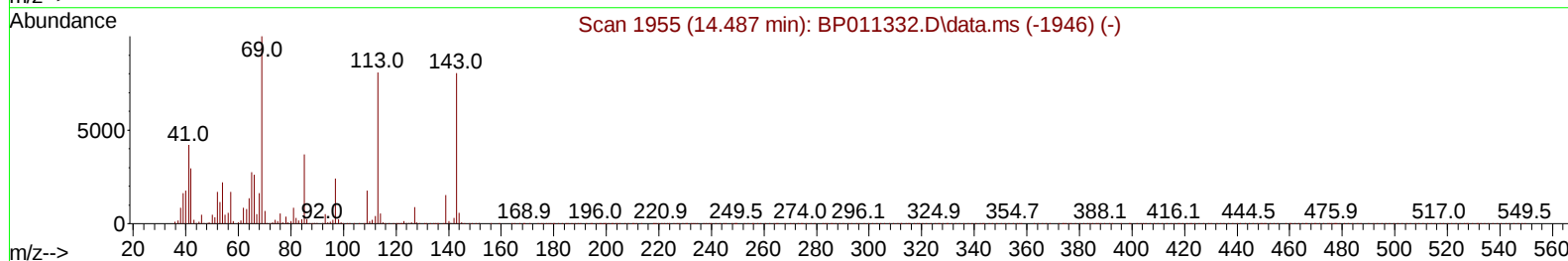
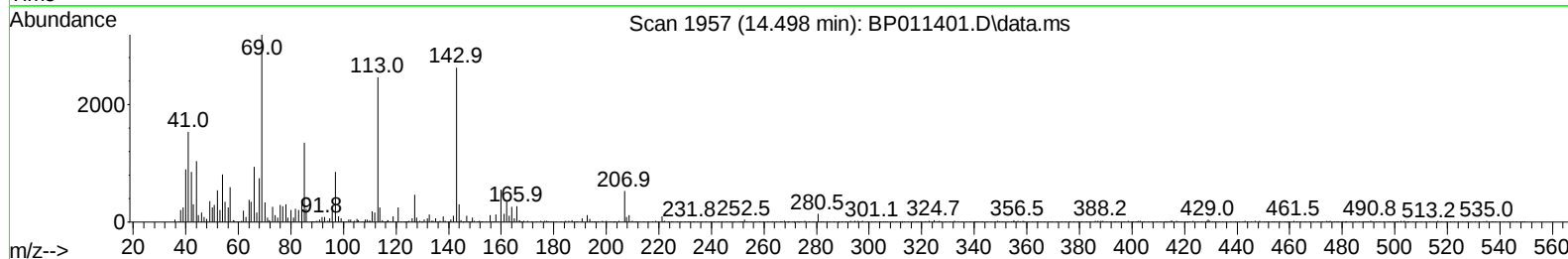
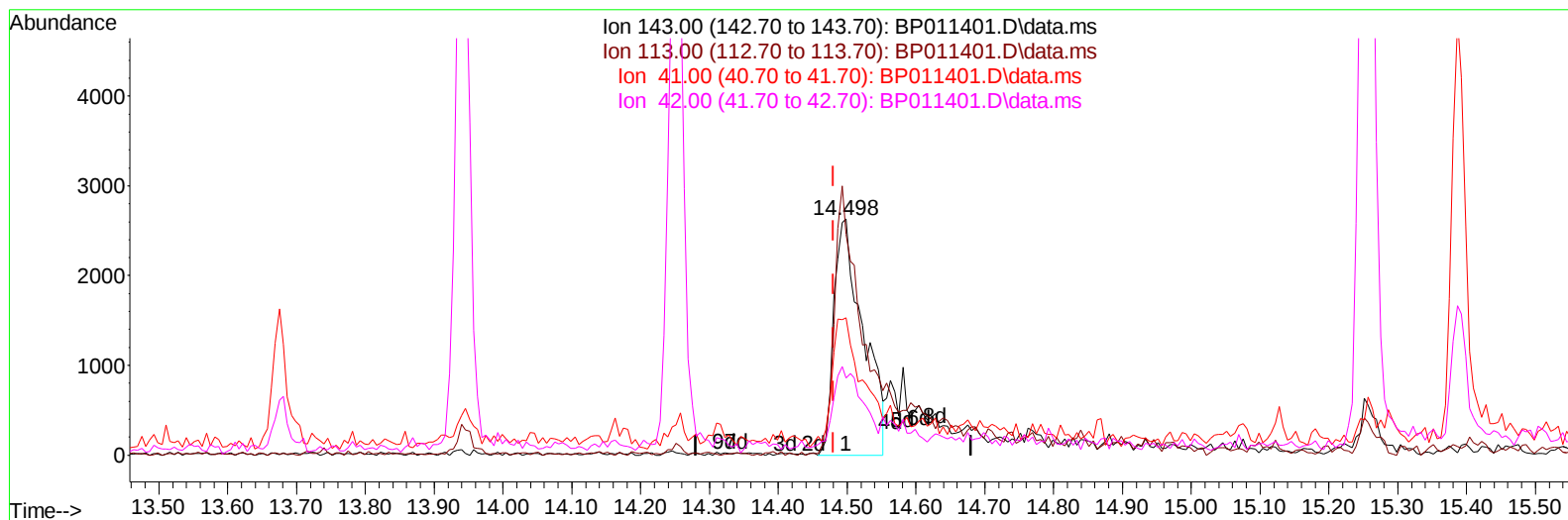
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TIC: BP011401.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 3.00 ng/ul m

response 7737

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	93.50#
41.00	54.40	58.19
42.00	35.50	32.57

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	84129	20.000	ng/ul	0.00
20) Naphthalene-d8	10.369	136	352536	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.251	164	203408	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	448511	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.151	240	478408	20.000	ng/ul	# 0.00
88) Perylene-d12	23.457	264	518867	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	11710	4.267	ng/uL	0.00
4) Pyridine-d5	3.499	84	48749	7.042	ng/ul	0.01
7) Phenol-d5	6.764	99	43257	5.810	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	123853	23.461	ng/ul	0.00
11) 2-Chlorophenol-d4	7.111	132	125084	21.412	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	77671	13.287	ng/ul	0.00
21) Nitrobenzene-d5	8.746	128	70804	29.511	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	67279	30.628	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.999	165	108586	24.478	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	137062	17.673	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	456591	32.352	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	507941	29.898	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	7737m	3.001	ng/ul	0.02
60) Fluorene-d10	15.257	176	373263	30.925	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	56942	27.333	ng/ul	0.00
73) Anthracene-d10	17.122	188	705100	34.976	ng/ul	0.00
81) Pyrene-d10	19.386	212	848843	32.205	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	914967	34.967	ng/ul	0.00

Target Compounds Qvalue

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

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