

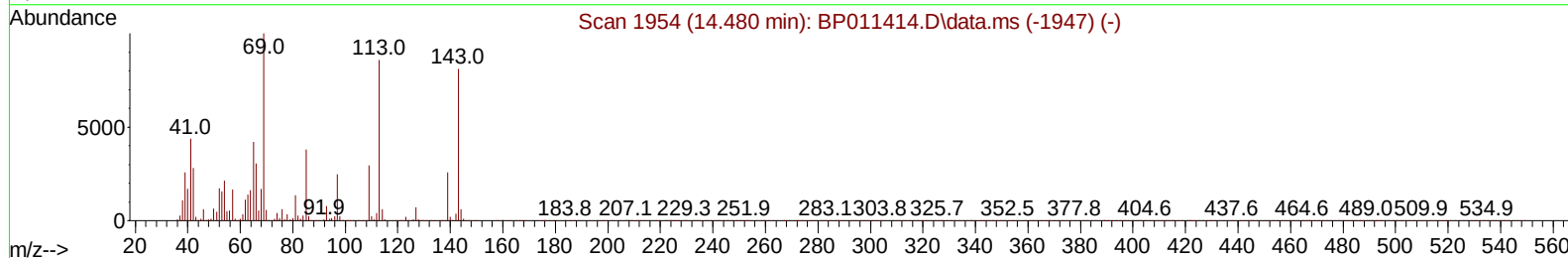
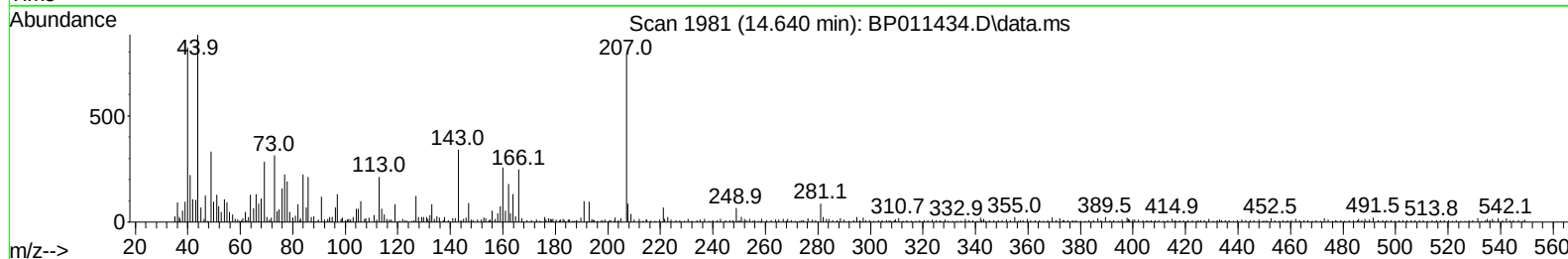
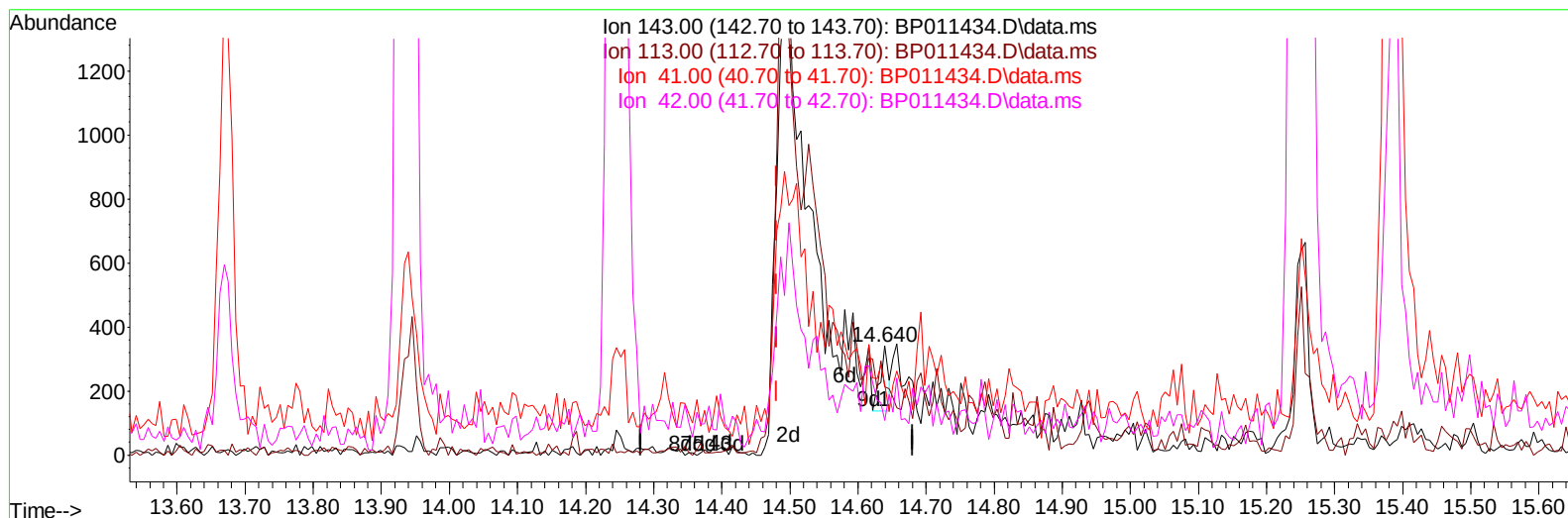
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011434.D
 Acq On : 12 Aug 2022 20:59
 Operator : CG/JU
 Sample : N4165-17
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBY9

Manual IntegrationsAPPROVED

Quant Time: Aug 12 23:18:25 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 :Yogesh Patel 08/13/2022
 Supervised By :Sohil Jodhani 08/13/2022



TIC: BP011434.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.640min (+ 0.159) 0.07 ng/uI

response 164

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	62.46
41.00	54.40	64.52
42.00	35.50	31.96

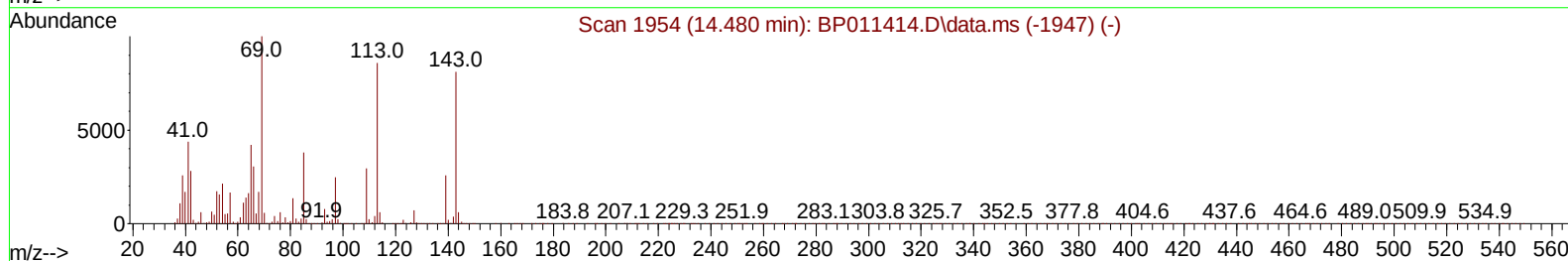
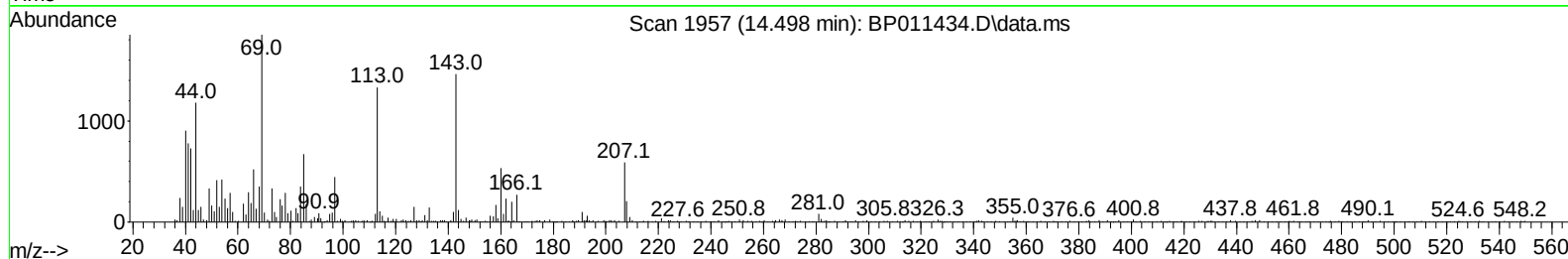
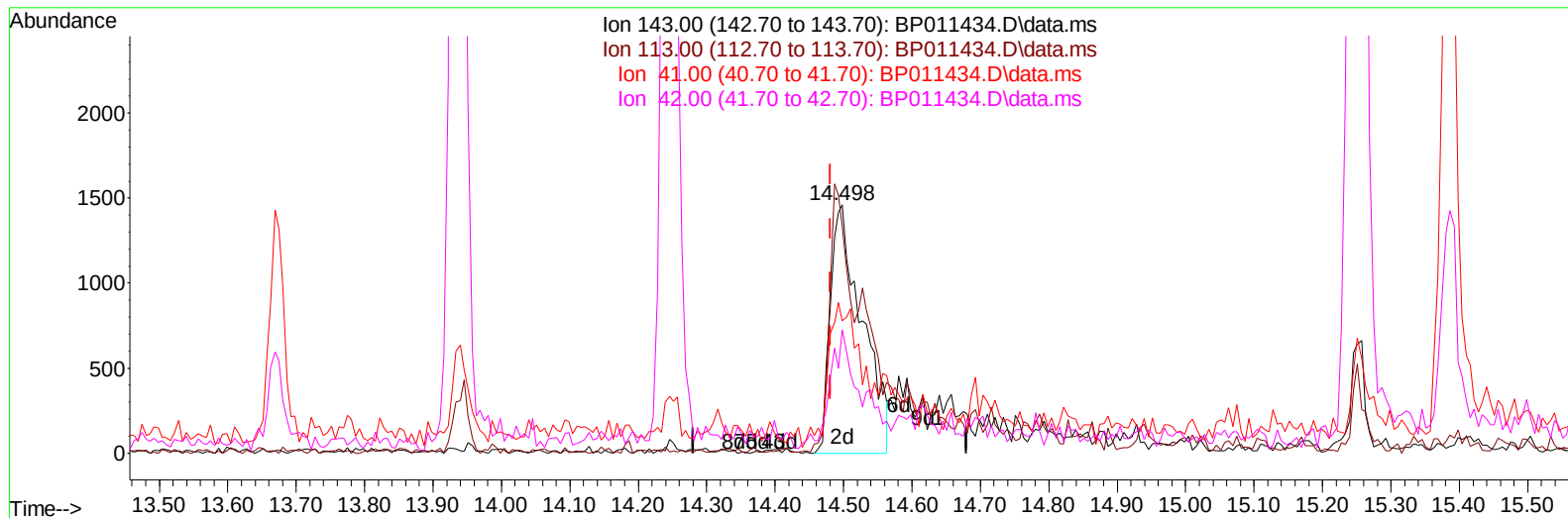
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(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 2.12 ng/ul m

response 4708

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	91.23#
41.00	54.40	53.36
42.00	35.50	49.66#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	70181	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	304025	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.245	164	175489	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	362163	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.151	240	351525	20.000	ng/ul	0.00
88) Perylene-d12	23.457	264	367763	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	11128	4.861	ng/uL	0.00
4) Pyridine-d5	3.499	84	40143	6.951	ng/ul	0.01
7) Phenol-d5	6.758	99	30813	4.961	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	127938	29.052	ng/ul	0.00
11) 2-Chlorophenol-d4	7.105	132	105993	21.750	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	60350	12.375	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	70452	34.049	ng/ul	0.00
24) 2-Nitrophenol-d4	9.457	143	64584	34.093	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	97593	25.510	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	152007	22.728	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	435592	35.774	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	453110	30.914	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	4708m	2.117	ng/ul	0.02
60) Fluorene-d10	15.251	176	357023	34.286	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	48033	28.554	ng/ul	0.00
73) Anthracene-d10	17.122	188	620710	38.131	ng/ul	0.00
81) Pyrene-d10	19.380	212	711926	36.760	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.304	264	703917	37.954	ng/ul	-0.01

Target Compounds Qvalue

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

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