

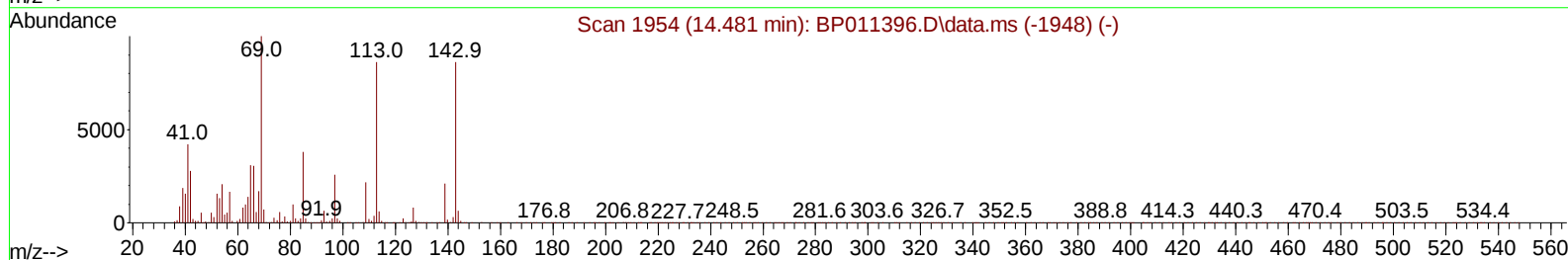
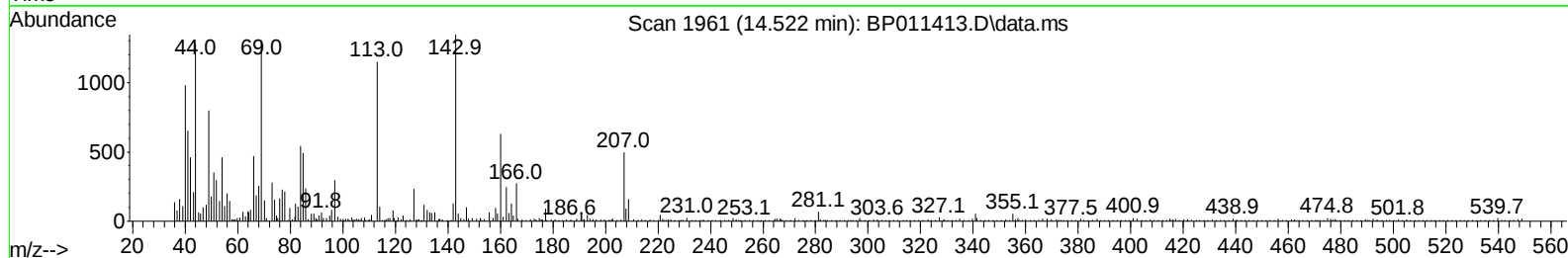
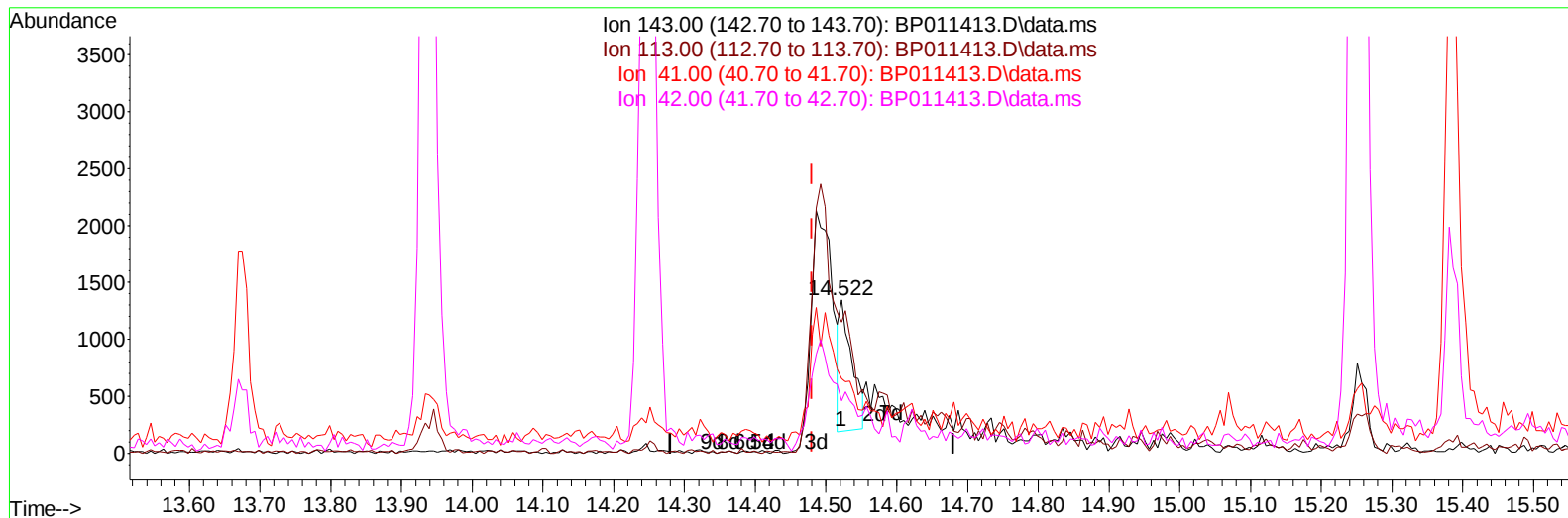
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011413.D
 Acq On : 12 Aug 2022 06:33
 Operator : CG/JU
 Sample : N4129-05
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBS8

Manual IntegrationsAPPROVED

Quant Time: Aug 12 07:18:48 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: BP011413.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.041) 0.60 ng/u1

response 1401

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	85.60
41.00	54.40	48.63
42.00	35.50	34.30

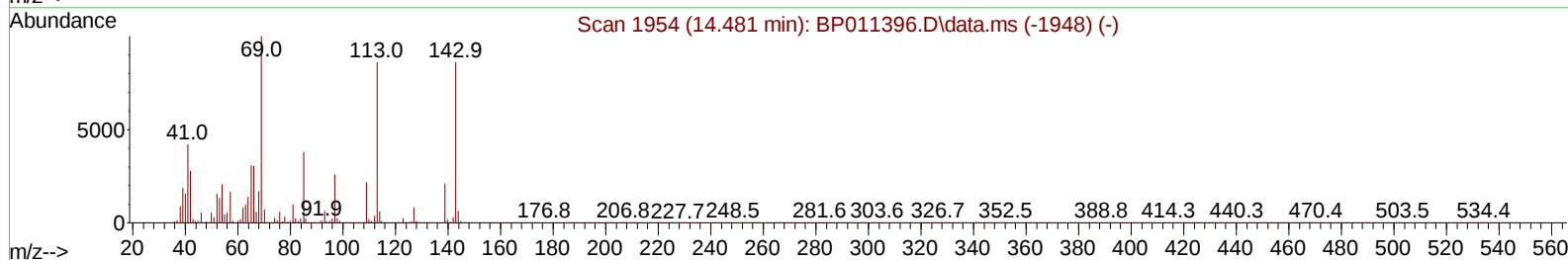
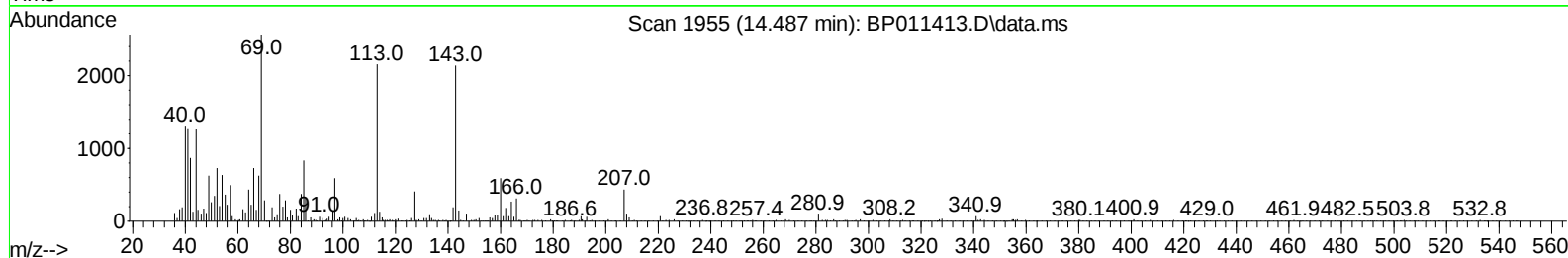
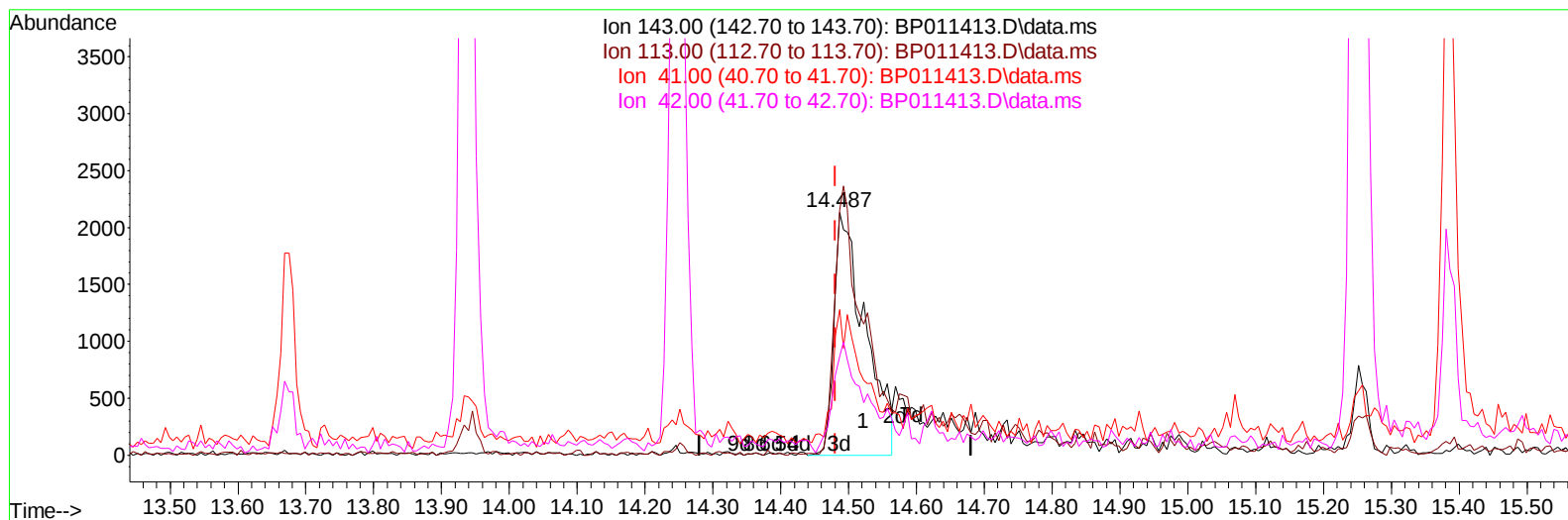
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TIC: BP011413.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.006) 2.85 ng/ul m

response 6705

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	100.84#
41.00	54.40	60.03
42.00	35.50	40.53

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	76853	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	324995	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	185333	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	382698	20.000	ng/ul	0.00
79) Chrysene-d12	21.151	240	374272	20.000	ng/ul	0.00
88) Perylene-d12	23.457	264	386956	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	10848	4.327	ng/uL	0.00
4) Pyridine-d5	3.499	84	53243	8.420	ng/ul	0.01
7) Phenol-d5	6.758	99	41802	6.146	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	147702	30.628	ng/ul	0.00
11) 2-Chlorophenol-d4	7.111	132	132744	24.875	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	79165	14.824	ng/ul	0.00
21) Nitrobenzene-d5	8.740	128	82538	37.317	ng/ul	0.00
24) 2-Nitrophenol-d4	9.458	143	79096	39.059	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	121902	29.809	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	162868	22.780	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	477210	37.110	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	554483	35.821	ng/ul	0.00
54) 4-Nitrophenol-d4	14.487	143	6705m	2.855	ng/ul	0.00
60) Fluorene-d10	15.251	176	393582	35.789	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	52662	29.626	ng/ul	0.00
73) Anthracene-d10	17.122	188	667229	38.789	ng/ul	0.00
81) Pyrene-d10	19.386	212	771873	37.433	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.310	264	755209	38.700	ng/ul	0.00

Target Compounds Qvalue

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

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