

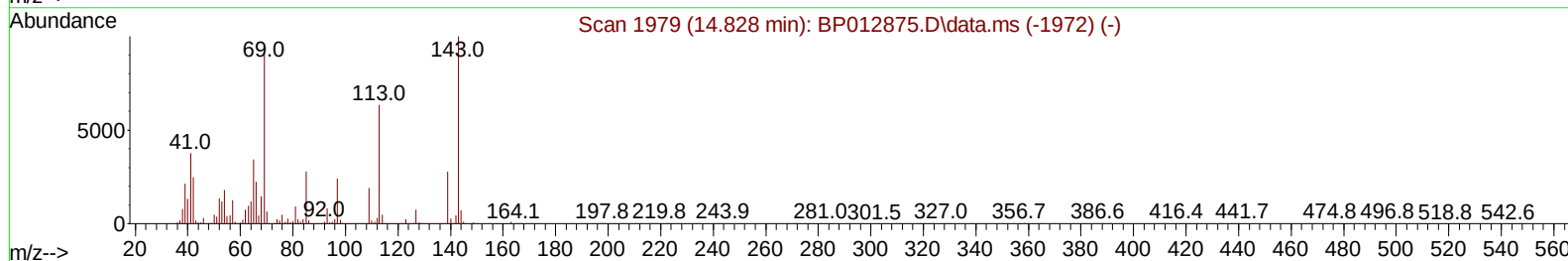
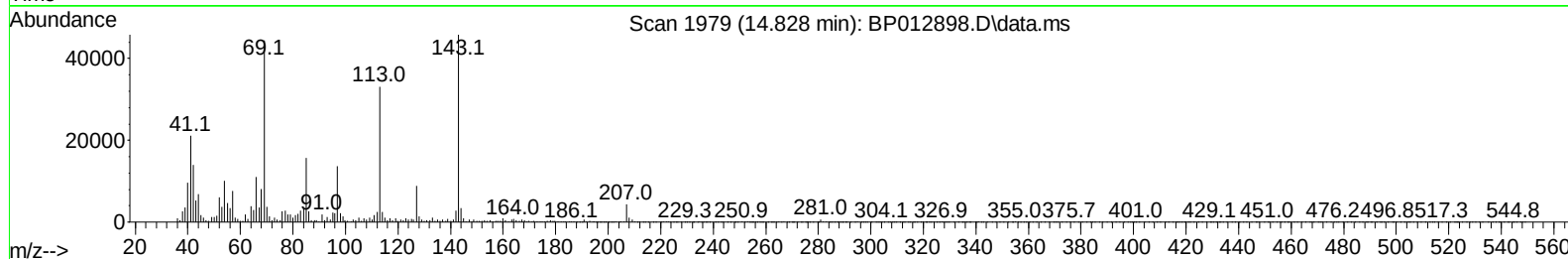
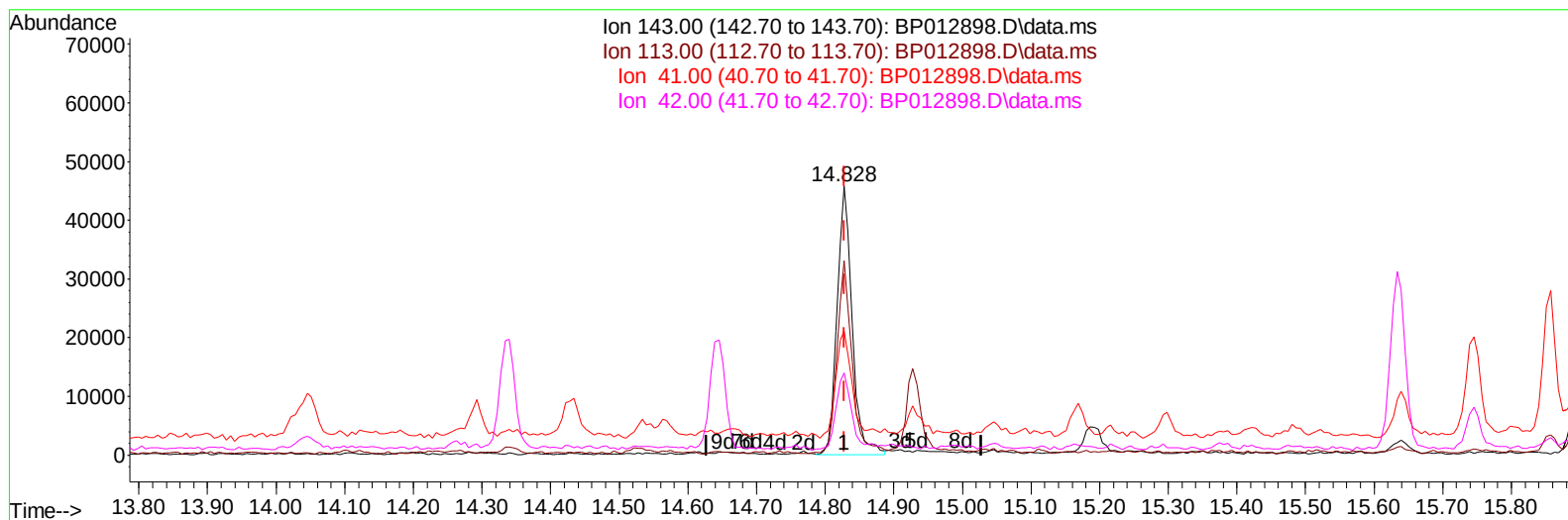
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012898.D
 Acq On : 01 Dec 2022 02:21
 Operator : CG/JU
 Sample : N5737-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4347

Manual IntegrationsAPPROVED

Quant Time: Dec 01 02:59:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/01/2022
 Supervised By :mohammad ahmed 12/02/2022



TIC: BP012898.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.828min (-0.000) 5.01 ng/ul m

response 69802

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.70	72.24
41.00	37.00	46.06#
42.00	24.70	30.56#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	456831	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	1835255	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1136398	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	2440706	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	2504645	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	2618798	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	52413	4.916	ng/uL	0.00
4) Pyridine-d5	3.822	84	367973	11.534	ng/ul	0.00
7) Phenol-d5	7.157	99	251228	6.692	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	644512	27.863	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	662730	23.169	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	442969	14.646	ng/ul	0.00
21) Nitrobenzene-d5	9.175	128	396186	35.270	ng/ul	0.00
24) 2-Nitrophenol-d4	9.899	143	395269	33.140	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	748794	26.907	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	915254	24.017	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	2499159	31.950	ng/ul	0.00
49) Acenaphthylene-d8	14.339	160	2821380	30.913	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	69802m	5.005	ng/ul	0.00
60) Fluorene-d10	15.634	176	2220051	32.408	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	288614	24.741	ng/ul	0.00
73) Anthracene-d10	17.498	188	3547977	33.024	ng/ul	0.00
81) Pyrene-d10	19.727	212	4211613	28.939	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	4252539	33.059	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.428	88	549486	47.993	ng/uL	98

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

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