

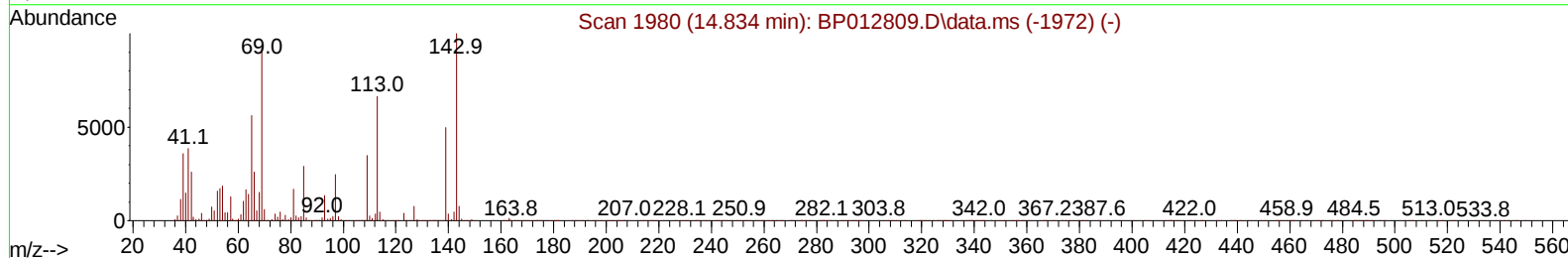
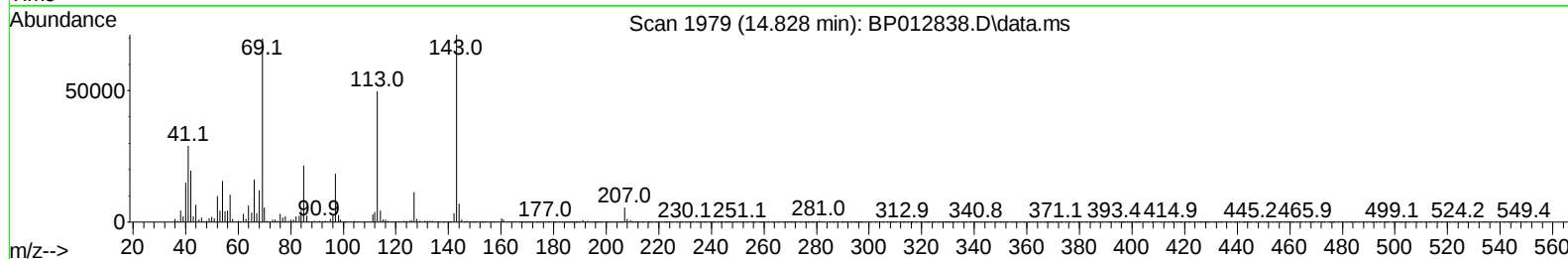
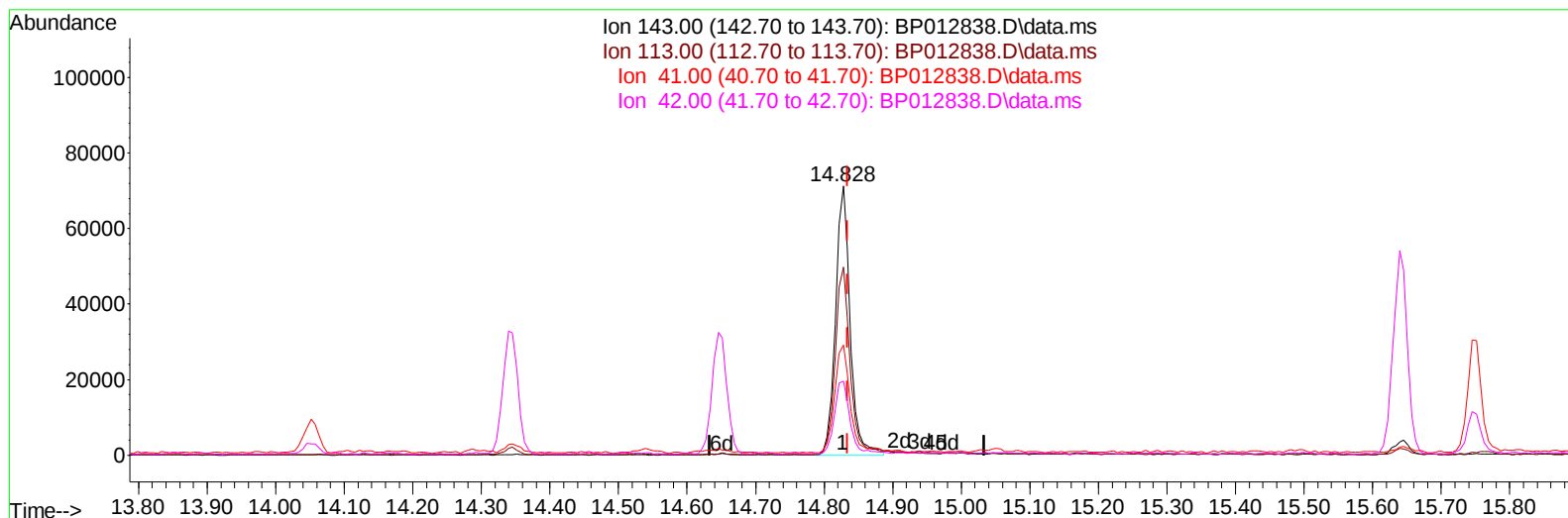
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
 Data File : BP012838.D
 Acq On : 29 Nov 2022 03:54
 Operator : CG/JU
 Sample : N5666-08
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB27

Manual IntegrationsAPPROVED

Quant Time: Nov 29 04:43:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/29/2022
 Supervised By :mohammad ahmed 11/30/2022



TIC: BP012838.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.828min (-0.006) 4.60 ng/ul m

response 105779

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.70	69.81
41.00	37.00	40.81
42.00	24.70	27.59

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	612957	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2774109	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1872157	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	3903860	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	2656354	20.000	ng/ul	-0.01
88) Perylene-d12	23.998	264	2146914	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	57677	4.032	ng/uL	0.00
4) Pyridine-d5	3.822	84	289240	6.757	ng/ul	0.00
7) Phenol-d5	7.163	99	387467	7.693	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	981701	31.630	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	955885	24.906	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	684532	16.868	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	589953	34.746	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	628031	34.835	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	1153879	27.430	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	1130300	19.622	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	4112831	31.916	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	4711736	31.337	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	105779m	4.604	ng/ul	0.00
60) Fluorene-d10	15.639	176	3710611	32.879	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	494417	26.498	ng/ul	0.00
73) Anthracene-d10	17.504	188	5791450	33.702	ng/ul	0.00
81) Pyrene-d10	19.727	212	6006728	38.917	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	3563005	33.787	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.434	88	100787	6.561	ng/uL	99

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

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