

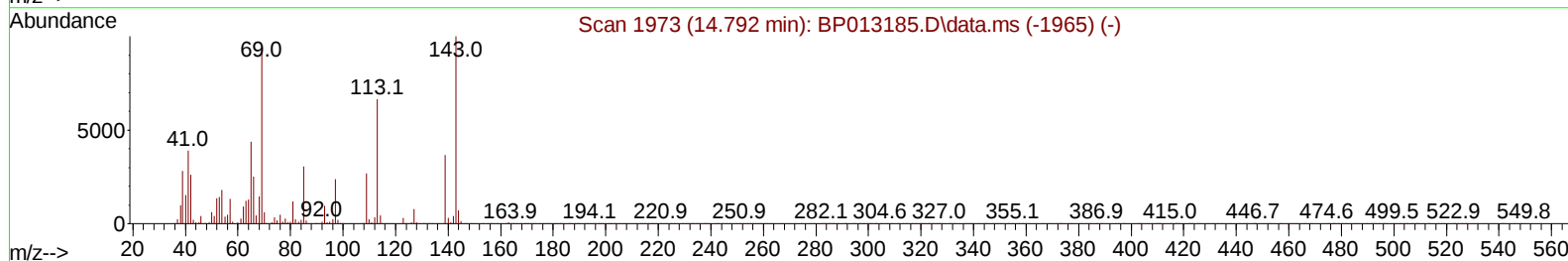
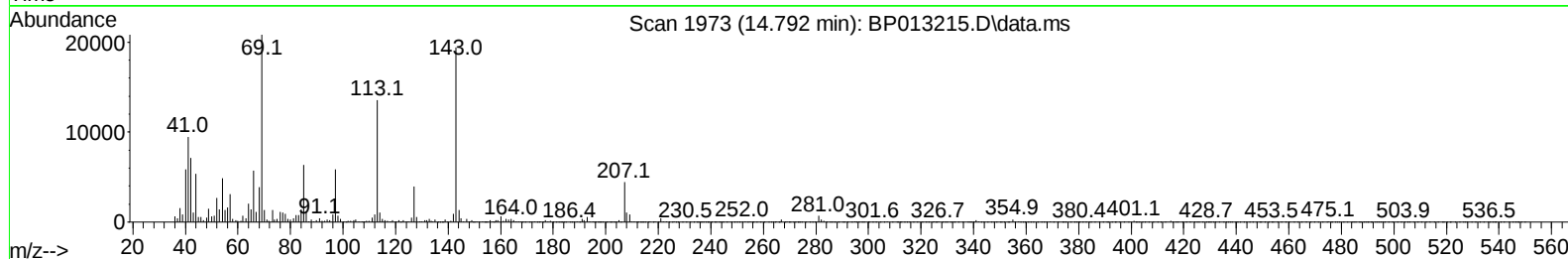
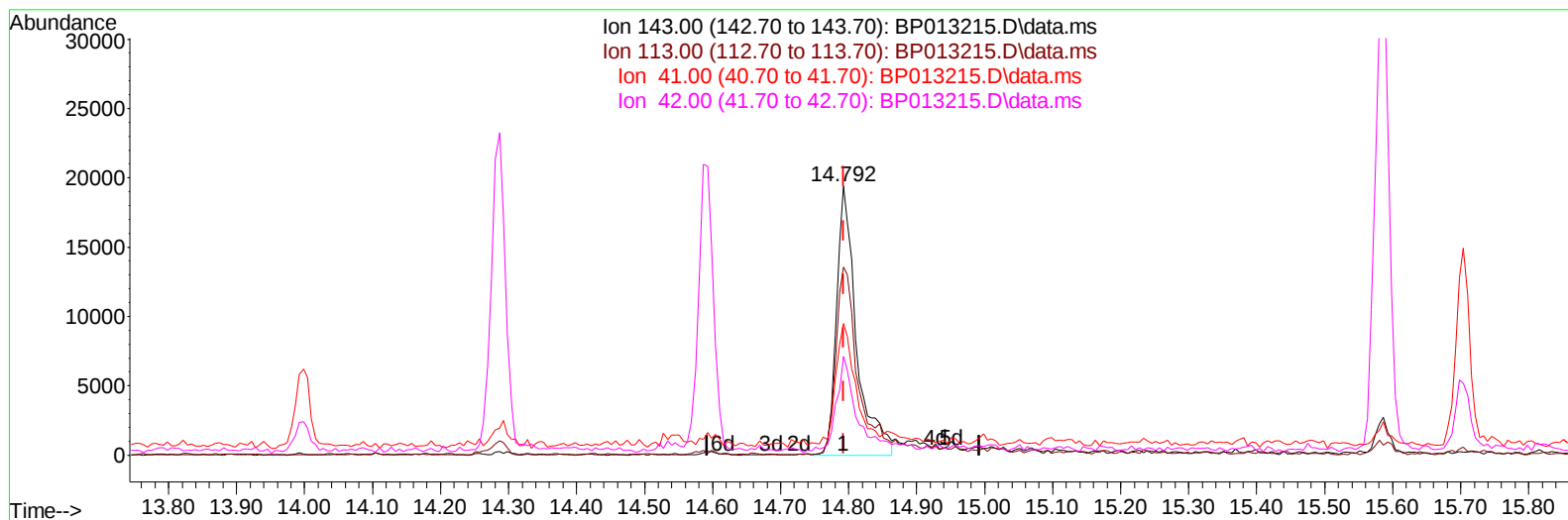
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013215.D
 Acq On : 21 Dec 2022 19:12
 Operator : CG/JU
 Sample : N6032-02
 Misc :
 ALS Vial : 97 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COL85

Manual IntegrationsAPPROVED

Quant Time: Dec 21 21:54:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 19 21:43:11 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/22/2022
 Supervised By :mohammad ahmed 12/22/2022



TIC: BP013215.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.792min (0.000) 2.69 ng/ul m

response 39346

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.80	69.89
41.00	35.70	48.85#
42.00	24.70	36.70#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	477746	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	1900384	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	1059846	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	2002245	20.000	ng/ul	-0.01
79) Chrysene-d12	21.433	240	2004235	20.000	ng/ul	0.00
88) Perylene-d12	23.910	264	2381241	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	1336710	119.631	ng/uL	-0.04
4) Pyridine-d5	3.781	84	45119	1.421	ng/ul	0.02
7) Phenol-d5	7.110	99	227191	5.838	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	649839	27.683	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	532477	17.382	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	340814	10.689	ng/ul	-0.01
21) Nitrobenzene-d5	9.122	128	427162	30.593	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	405460	25.794	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.381	165	638265	20.905	ng/ul	0.00
31) 4-Chloroaniline-d4	10.904	131	357536	8.295	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	2460100	30.202	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	2837345	31.701	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	39346m	2.685	ng/ul	0.00
60) Fluorene-d10	15.586	176	2131443	30.931	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	227908	17.688	ng/ul	0.00
73) Anthracene-d10	17.445	188	3138842	34.748	ng/ul	-0.01
81) Pyrene-d10	19.680	212	3710733	32.255	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	3957874	33.346	ng/ul	-0.01

Target Compounds Qvalue

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

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