

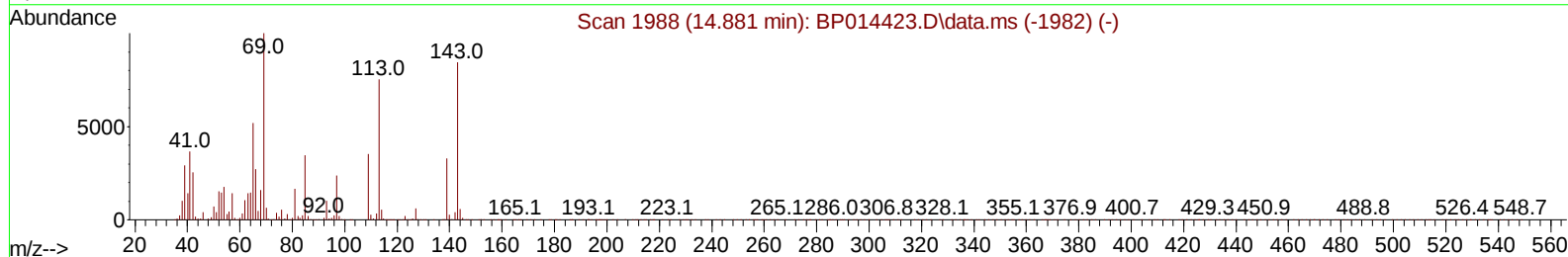
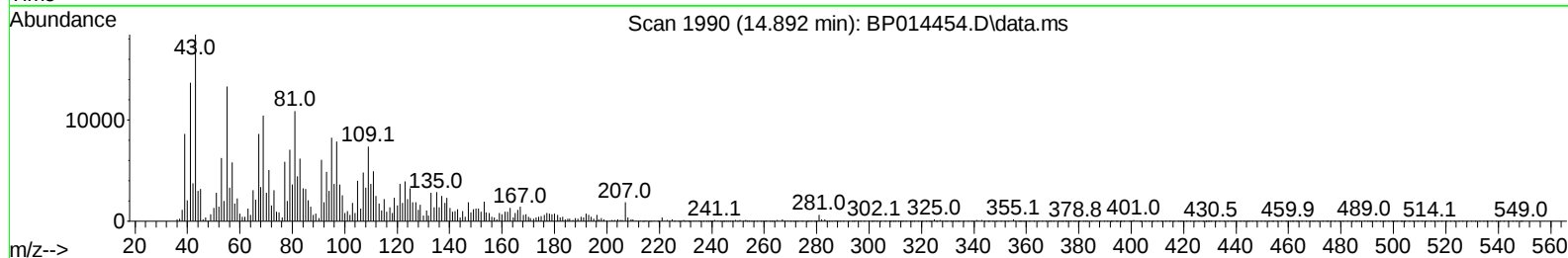
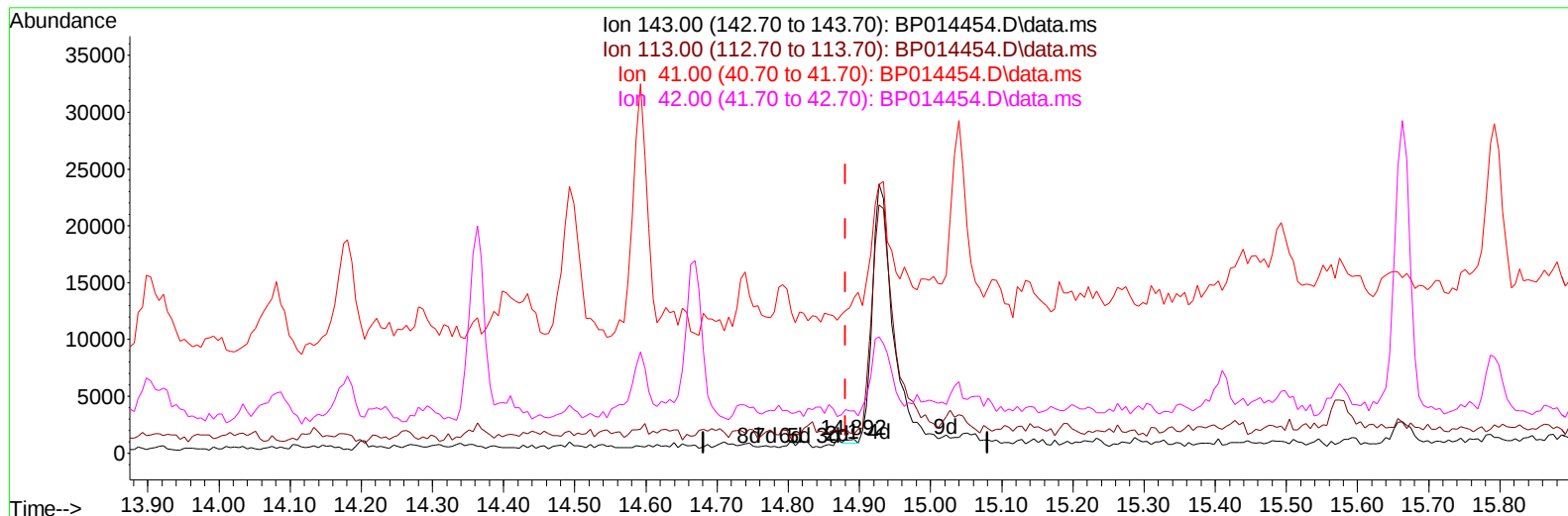
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014454.D
Acq On : 31 Mar 2023 21:03
Operator : CG/JU
Sample : 02092-09
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB973

Manual IntegrationsAPPROVED

Quant Time: Mar 31 22:56:13 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/03/2023
Supervised By :Jagrut Upadhyay 04/03/2023



TIC: BP014454.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.892min (+ 0.012) 0.04 ng/u1

response 307

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	143.31#
41.00	44.20	1138.57#
42.00	31.50	311.14#

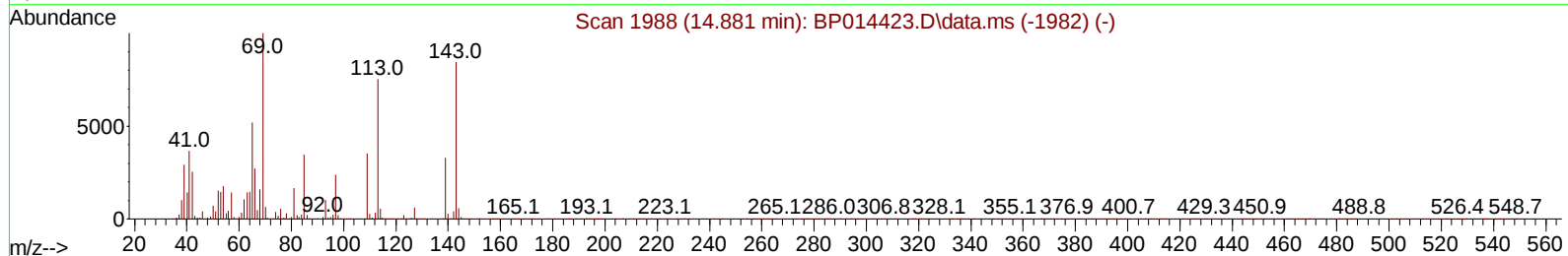
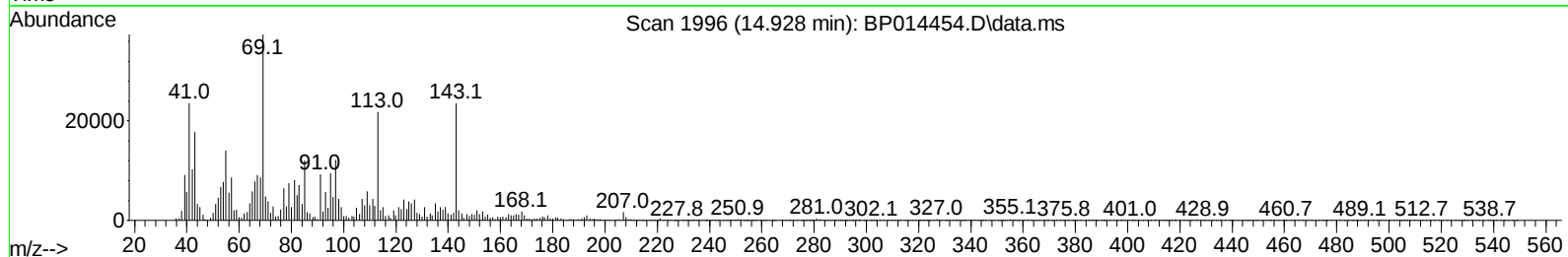
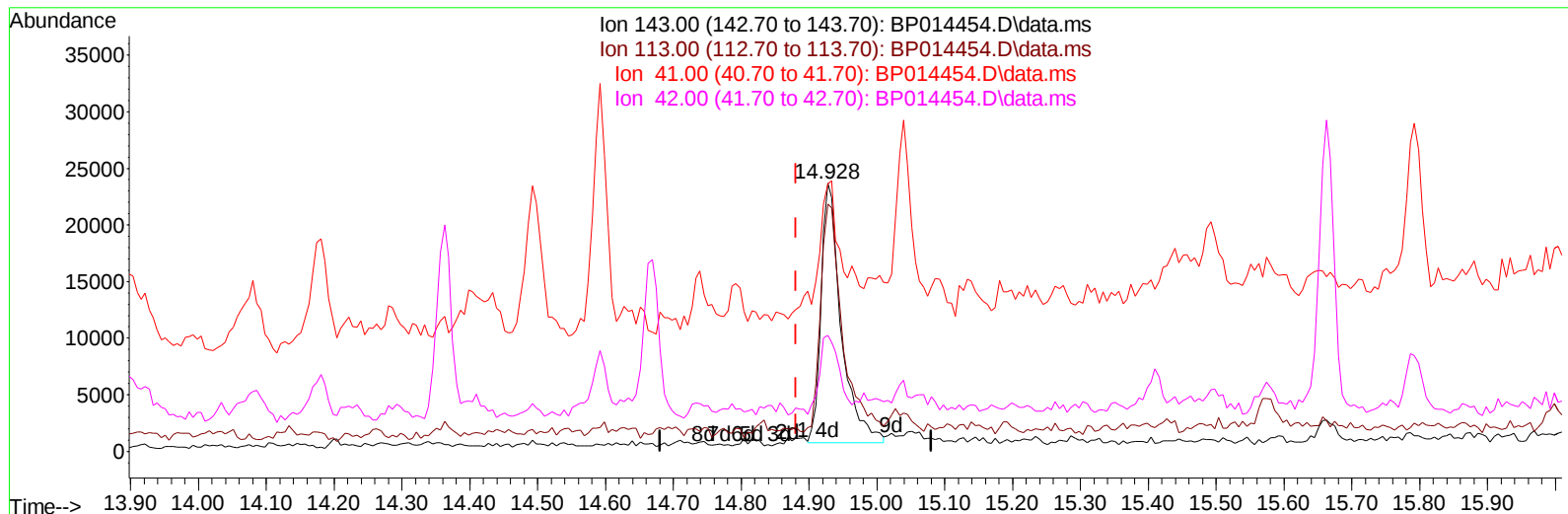
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(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.047) 6.28 ng/ul m

response 47440

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	92.19
41.00	44.20	99.93#
42.00	31.50	43.32#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	226782	20.000	ng/ul	0.00
20) Naphthalene-d8	10.840	136	977330	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	616830	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.433	188	1294220	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	876023	20.000	ng/ul	0.00
88) Perylene-d12	24.045	264	918978	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	22628	3.869	ng/uL	0.00
4) Pyridine-d5	3.828	84	145068	9.680	ng/ul	0.01
7) Phenol-d5	7.187	99	181225	8.665	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.340	67	395186	29.449	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	402376	26.618	ng/ul	0.00
15) 4-Methylphenol-d8	8.740	113	331052	19.623	ng/ul	0.01
21) Nitrobenzene-d5	9.193	128	249815	31.663	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	293205	32.454	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	486771	29.585	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	570228	26.311	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	1682828	33.647	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1823298	32.855	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	47440m	6.279	ng/ul	0.05
60) Fluorene-d10	15.663	176	1370647	32.863	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	251750	26.436	ng/ul	0.00
73) Anthracene-d10	17.533	188	2066552	32.994	ng/ul	0.00
81) Pyrene-d10	19.763	212	2114878	35.942	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.880	264	1618381	33.147	ng/ul	0.01

Target Compounds Qvalue

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

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