

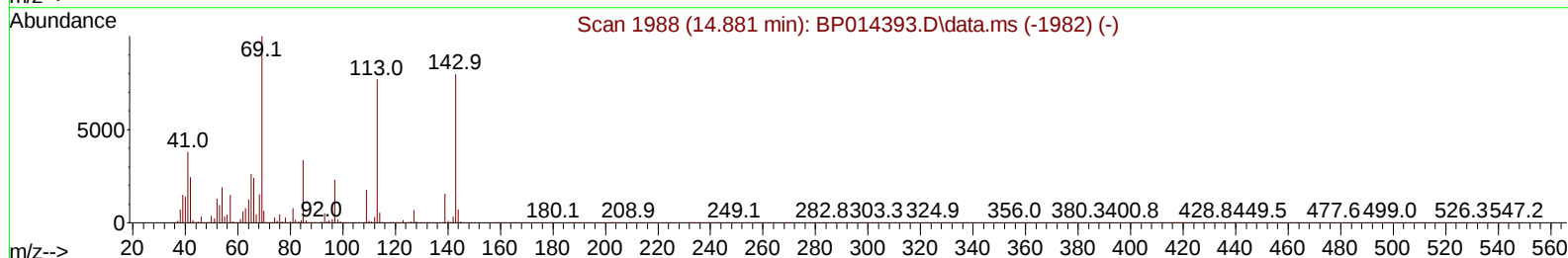
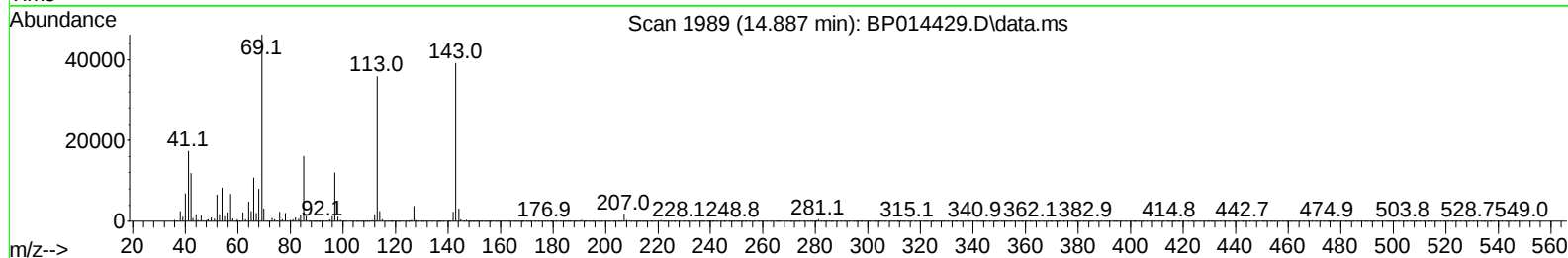
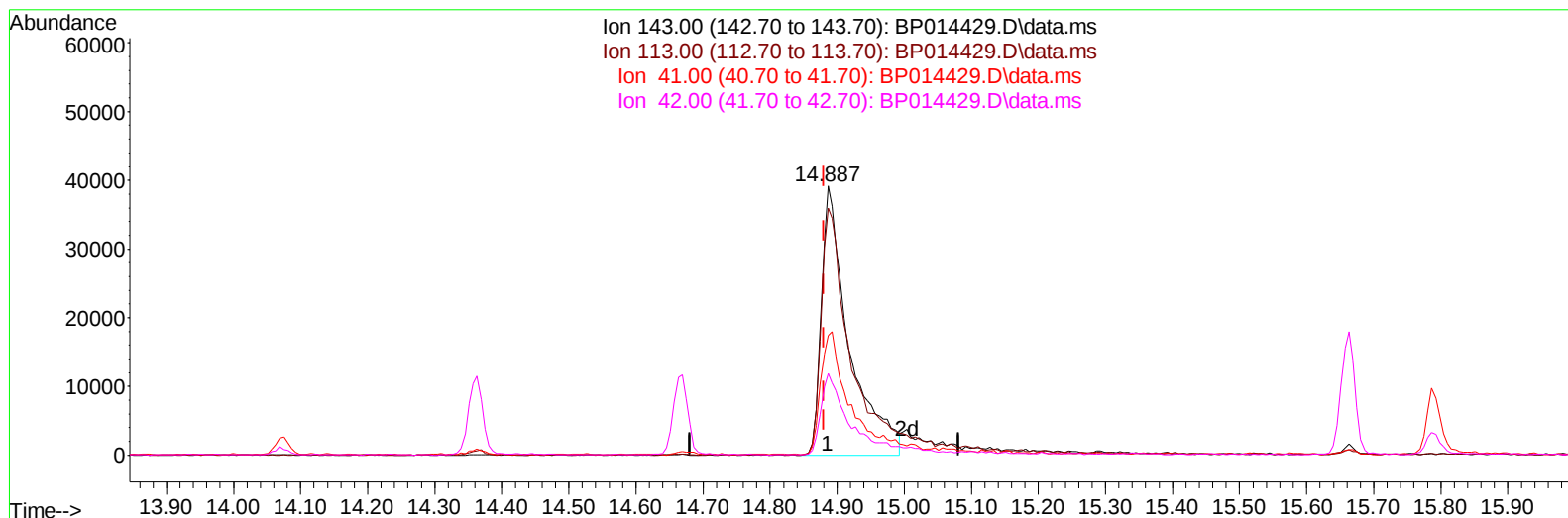
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014429.D
Acq On : 31 Mar 2023 06:58
Operator : CG/JU
Sample : PB151684BL
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK684

Manual IntegrationsAPPROVED

Quant Time: Mar 31 07:30:07 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 03/31/2023
Supervised By :Jagrut Upadhyay 03/31/2023



TIC: BP014429.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.006) 18.60 ng/u1

response 112142

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	91.85
41.00	44.20	44.44
42.00	31.50	30.31

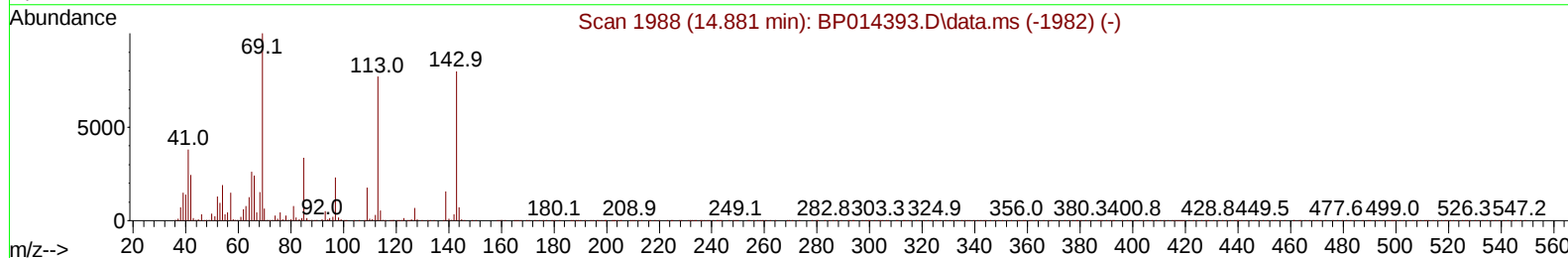
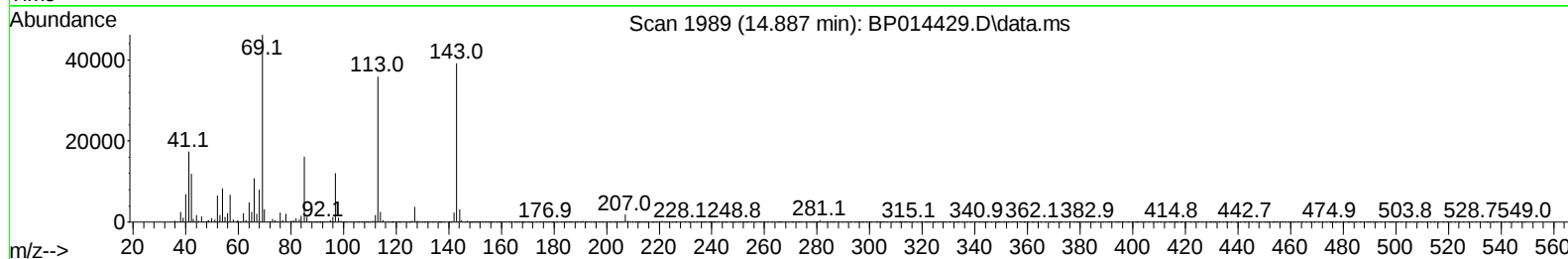
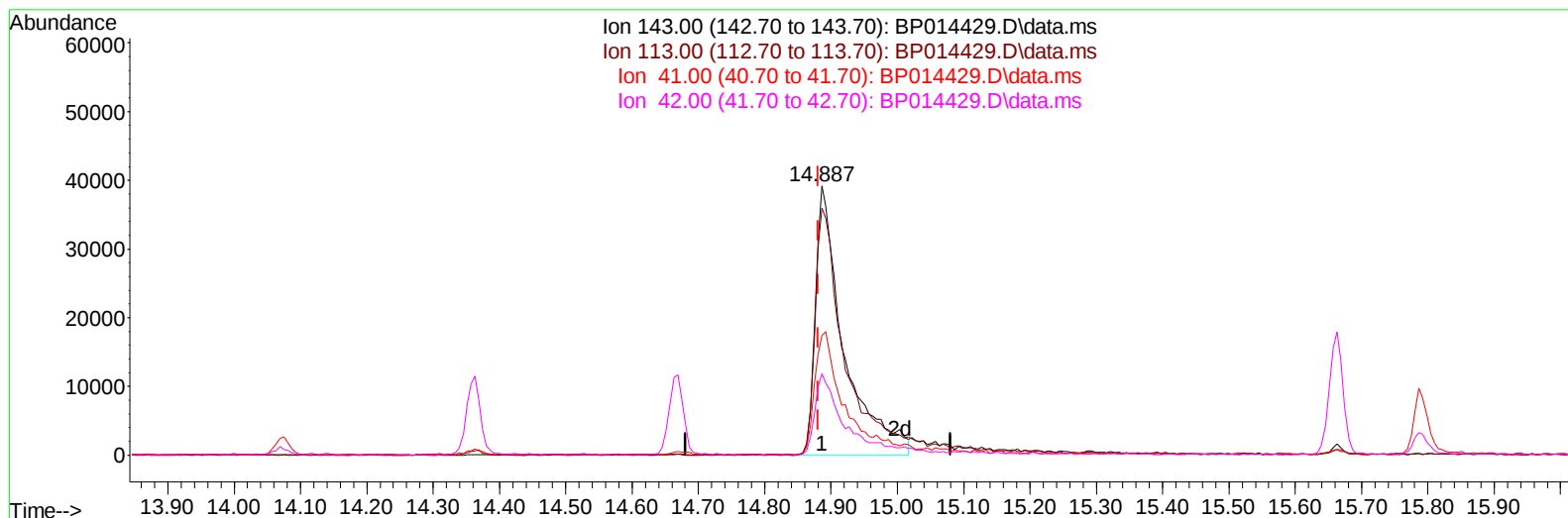
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TIC: BP014429.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.006) 19.30 ng/ul m

response 116380

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	91.85
41.00	44.20	44.44
42.00	31.50	30.31

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	186654	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	783235	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	492169	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	959814	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	813797	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	918610	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	25791	5.358	ng/uL	0.00
4) Pyridine-d5	3.816	84	328542	26.636	ng/ul	0.00
7) Phenol-d5	7.169	99	450058	26.145	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	307888	27.877	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	342667	27.542	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	356770	25.694	ng/ul	0.00
21) Nitrobenzene-d5	9.193	128	171876	27.183	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	204068	28.185	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	336756	25.539	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	447501	25.765	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	1091705	27.357	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1200023	27.101	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	116380m	19.304	ng/ul	0.00
60) Fluorene-d10	15.663	176	893767	26.857	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	163791	23.192	ng/ul	0.00
73) Anthracene-d10	17.528	188	1284499	27.653	ng/ul	0.00
81) Pyrene-d10	19.757	212	1407007	25.740	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	1371044	28.092	ng/ul	0.00

Target Compounds Qvalue

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

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