

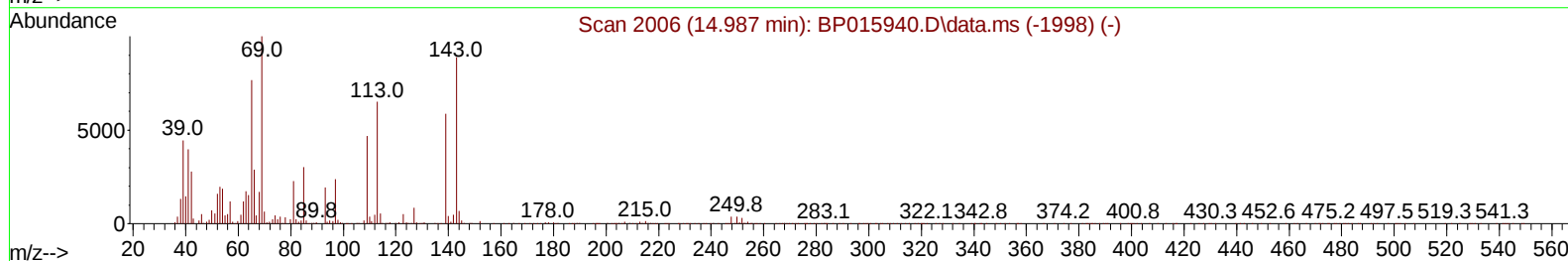
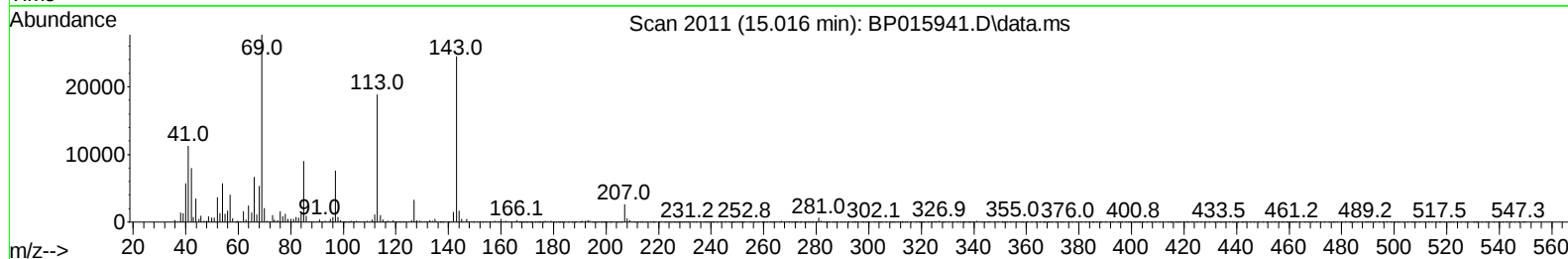
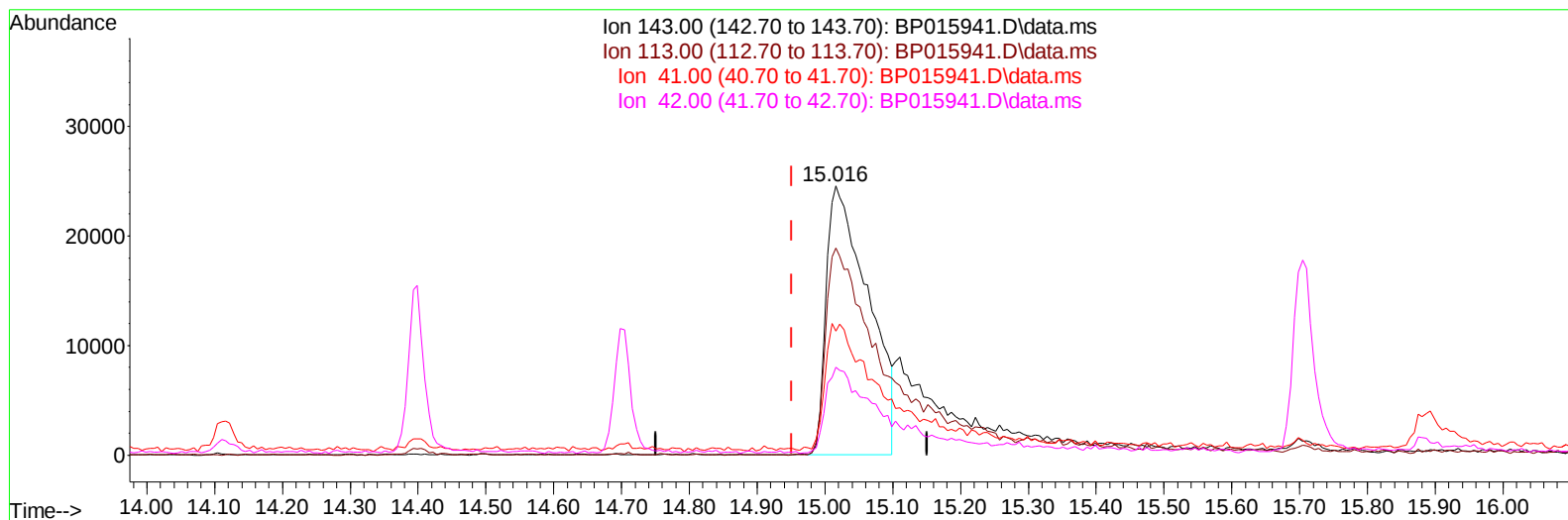
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070323\
 Data File : BP015941.D
 Acq On : 03 Jul 2023 09:23
 Operator : MA/JU
 Sample : PB153766BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK766

Manual IntegrationsAPPROVED

Quant Time: Jul 03 22:06:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jul 02 06:12:16 2023
 Response via : Initial Calibration

Reviewed By :Sohil Jodhani 07/04/2023
 Supervised By :Yogesh Patel 07/04/2023



TIC: BP015941.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.016min (+ 0.065) 19.53 ng/ul

response 105111

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	77.03
41.00	45.70	46.03
42.00	31.20	32.69

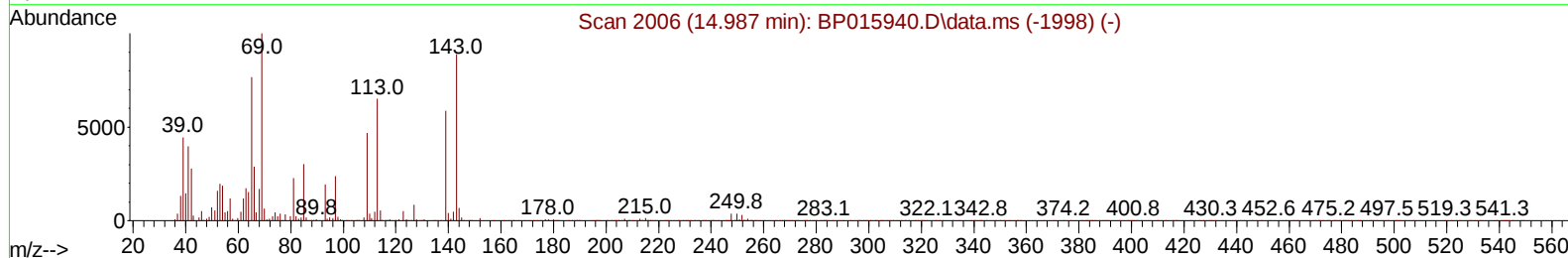
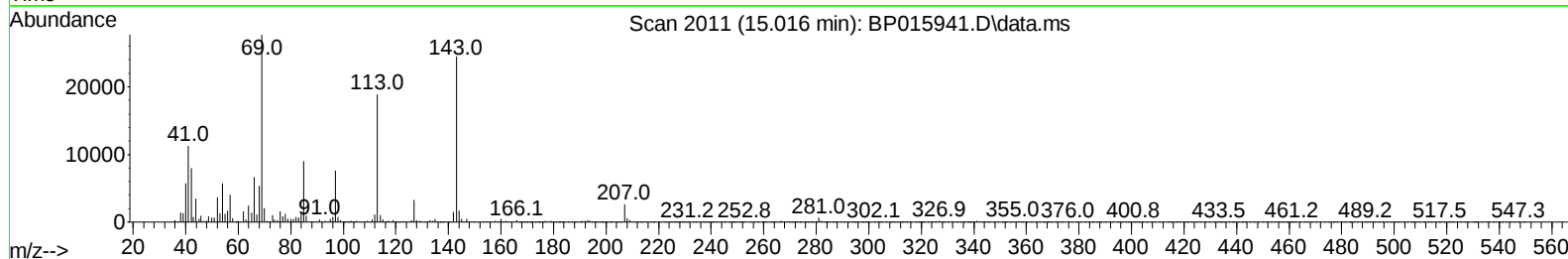
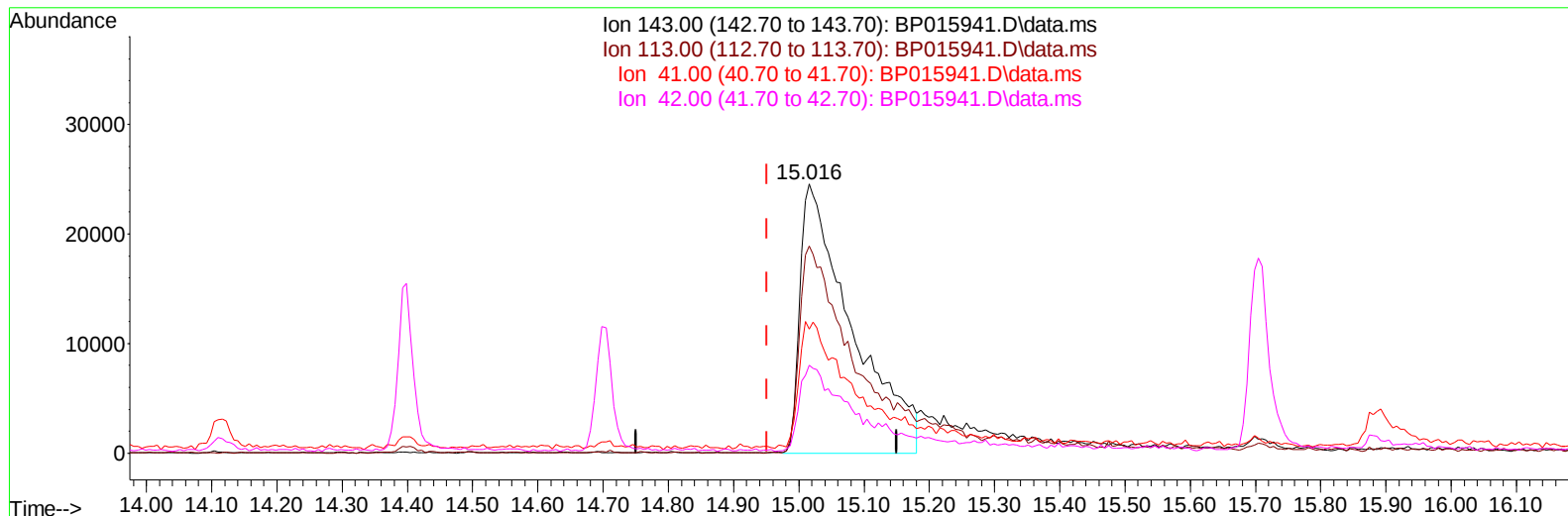
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TIC: BP015941.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.016min (+ 0.065) 25.06 ng/ul m

response 134858

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	77.03
41.00	45.70	46.03
42.00	31.20	32.69

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	196197	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136	903671	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.704	164	527642	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1003622	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.569	240	776368	20.000	ng/ul	0.02
88) Perylene-d12	24.110	264	847606	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	38523	7.064	ng/uL	0.00
4) Pyridine-d5	3.817	84	504513	36.702	ng/ul	0.00
7) Phenol-d5	7.234	99	619416	36.899	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.369	67	425022	40.924	ng/ul	0.00
11) 2-Chlorophenol-d4	7.581	132	511110	40.334	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	523590	39.482	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	244487	35.401	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143	261751	32.474	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.540	165	437835	30.482	ng/ul	0.03
31) 4-Chloroaniline-d4	11.052	131	648577	35.151	ng/ul	0.02
46) Dimethylphthalate-d6	14.116	166	1475345	36.176	ng/ul	0.01
49) Acenaphthylene-d8	14.398	160	1662503	36.263	ng/ul	0.00
54) 4-Nitrophenol-d4	15.016	143	134858m	25.058	ng/ul	0.07
60) Fluorene-d10	15.704	176	1198616	35.694	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.887	200	106697	17.287	ng/ul	0.03
73) Anthracene-d10	17.569	188	1771256	38.637	ng/ul	0.01
81) Pyrene-d10	19.798	212	1856544	36.622	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.939	264	1658197	38.782	ng/ul	0.01

Target Compounds Qvalue

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

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