

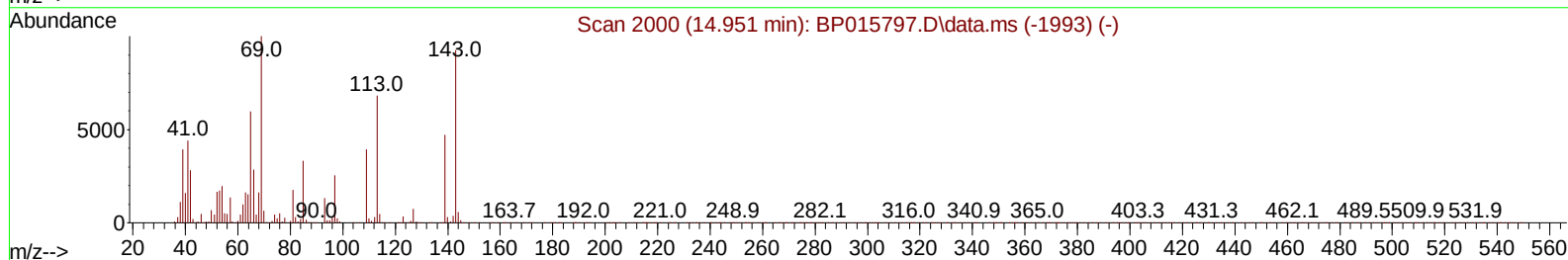
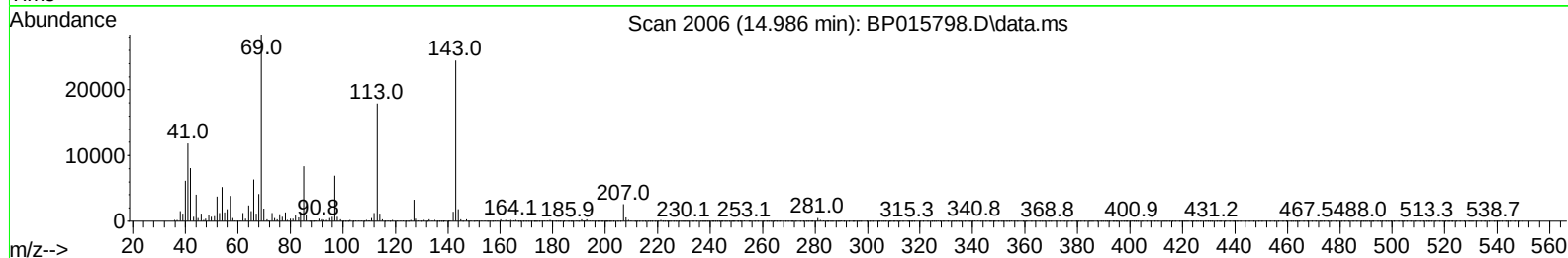
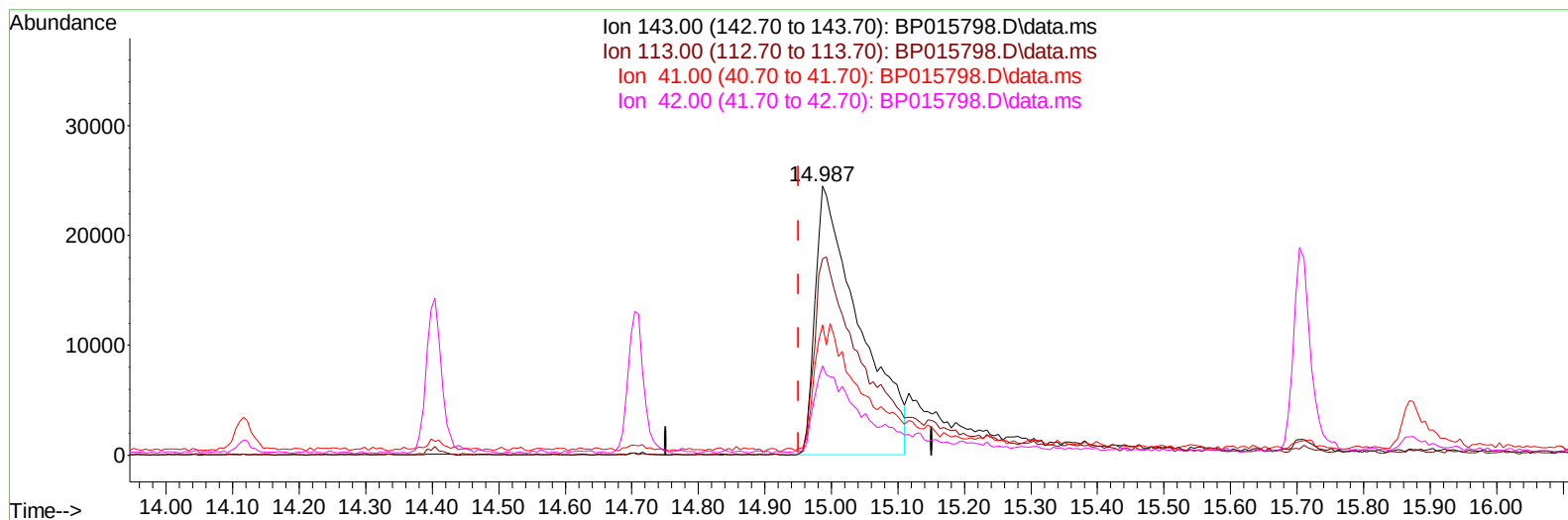
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015798.D
Acq On : 29 Jun 2023 07:52
Operator : MA/JU
Sample : PB153627BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK627

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/29/2023
Supervised By :mohammad ahmed 06/29/2023

Quant Time: Jun 29 08:27:22 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 29 02:30:23 2023
Response via : Initial Calibration



TIC: BP015798.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.986min (+ 0.035) 19.12 ng/ul

response 112842

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	73.09#
41.00	45.70	48.51
42.00	31.20	33.21

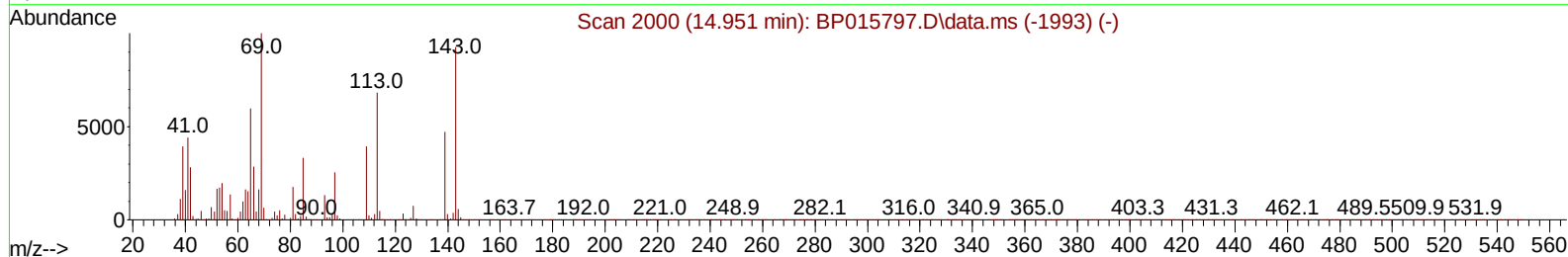
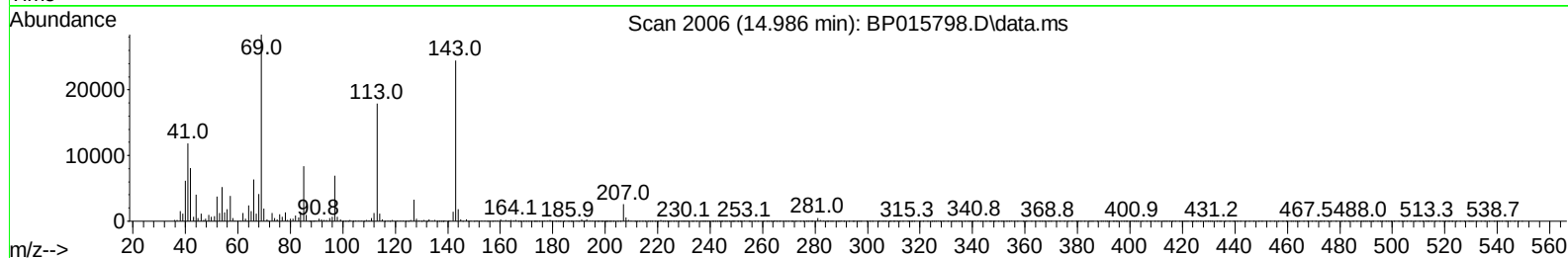
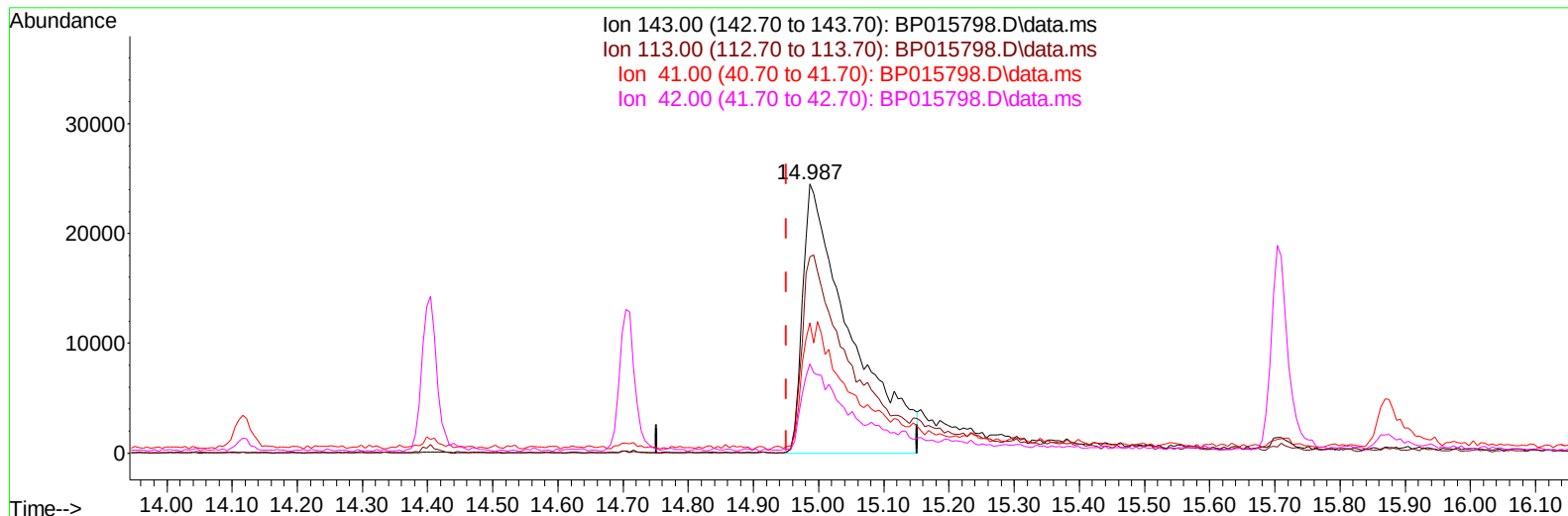
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TIC: BP015798.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.986min (+ 0.035) 21.02 ng/ul m

response 124073

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	73.09#
41.00	45.70	48.51
42.00	31.20	33.21

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Instrument :
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 QLast Update : Thu Jun 29 02:30:23 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.057	152	247814	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	1042416	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.704	164	578722	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	1040831	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.563	240	808411	20.000	ng/ul	# 0.00
88) Perylene-d12	24.104	264	943213	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	36650	5.321	ng/uL	0.00
4) Pyridine-d5	3.822	84	499733	28.782	ng/ul	0.00
7) Phenol-d5	7.228	99	635755	29.984	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67	439865	33.531	ng/ul	0.00
11) 2-Chlorophenol-d4	7.587	132	514335	32.134	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	525351	31.364	ng/ul	0.00
21) Nitrobenzene-d5	9.246	128	248602	31.206	ng/ul	0.00
24) 2-Nitrophenol-d4	9.975	143	270361	29.078	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.528	165	434087	26.199	ng/ul	0.01
31) 4-Chloroaniline-d4	11.051	131	638153	29.982	ng/ul	0.02
46) Dimethylphthalate-d6	14.116	166	1399722	31.292	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	1583695	31.495	ng/ul	0.00
54) 4-Nitrophenol-d4	14.986	143	124073m	21.019	ng/ul	0.04
60) Fluorene-d10	15.704	176	1114665	30.264	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	111239	17.378	ng/ul	0.02
73) Anthracene-d10	17.569	188	1559014	32.792	ng/ul	0.00
81) Pyrene-d10	19.798	212	1621452	30.717	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264	1614756	33.938	ng/ul	0.00

Target Compounds Qvalue

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

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