

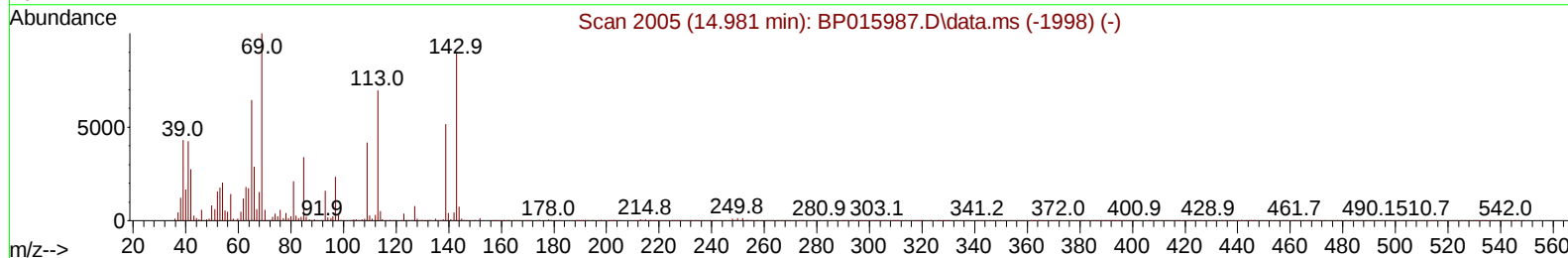
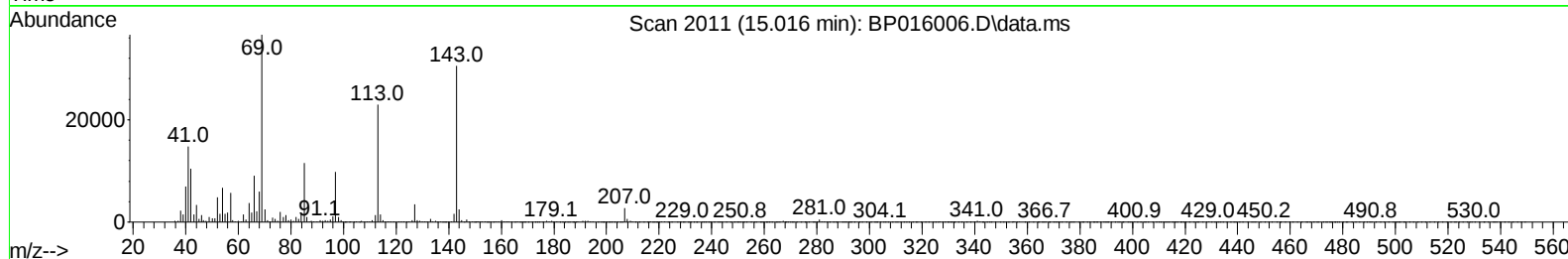
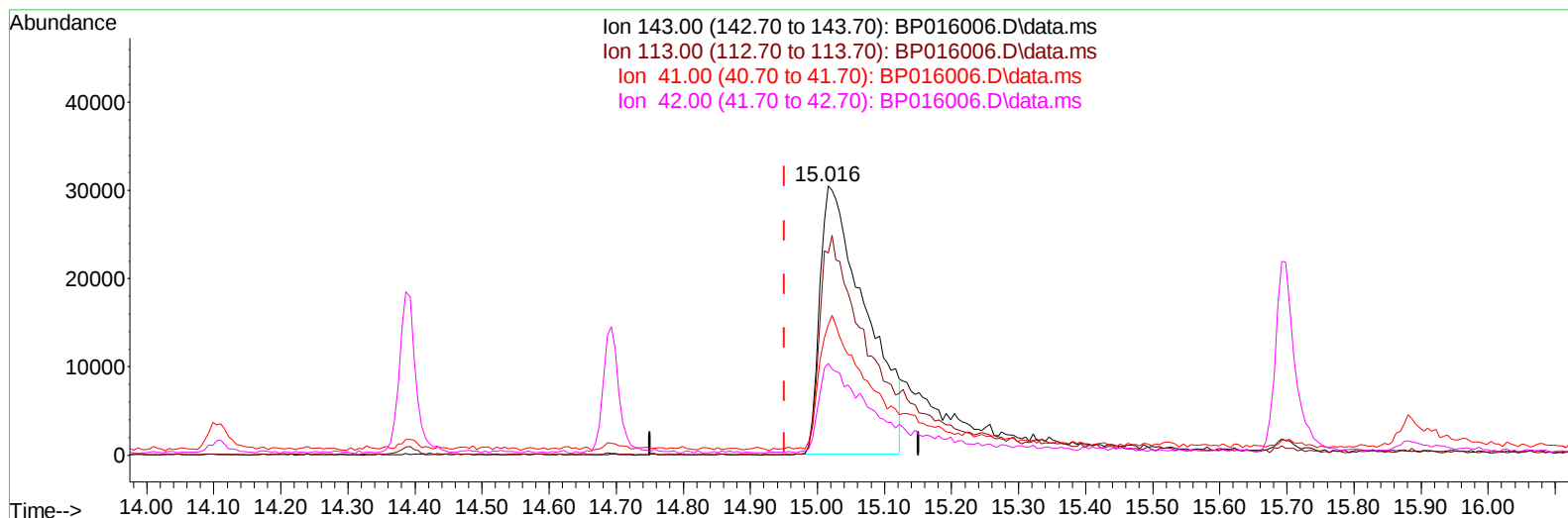
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP016006.D
Acq On : 05 Jul 2023 09:11
Operator : MA/JU
Sample : PB153469BL
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK469

Manual IntegrationsAPPROVED

Quant Time: Jul 05 11:33:35 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 :Yogesh Patel 07/05/2023
Supervised By :mohammad ahmed 07/05/2023



TIC: BP016006.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.016min (+ 0.065) 23.33 ng/u1

response 143345

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	75.18#
41.00	45.70	48.40
42.00	31.20	34.13

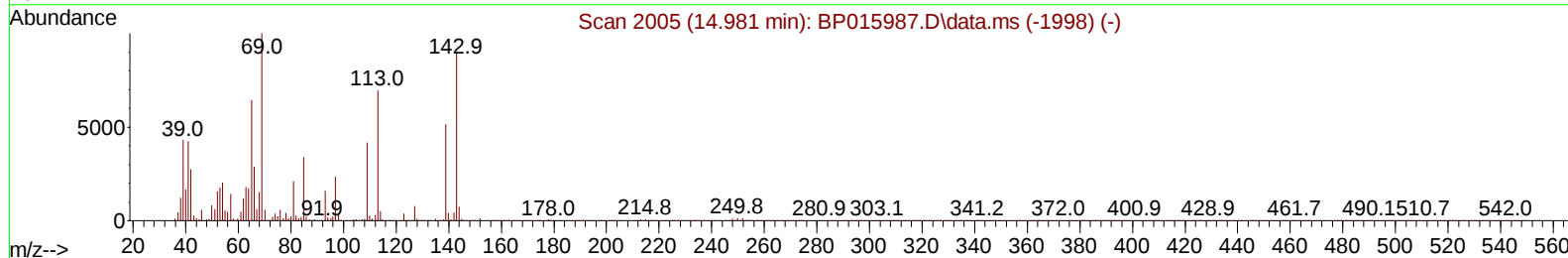
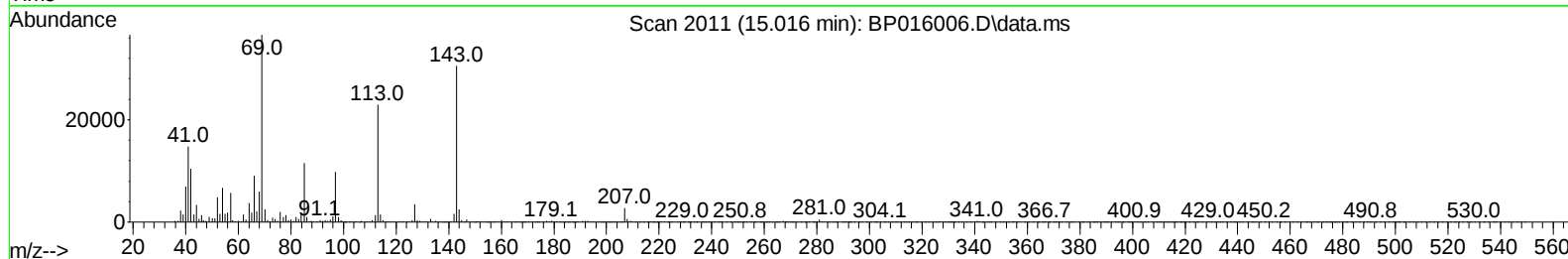
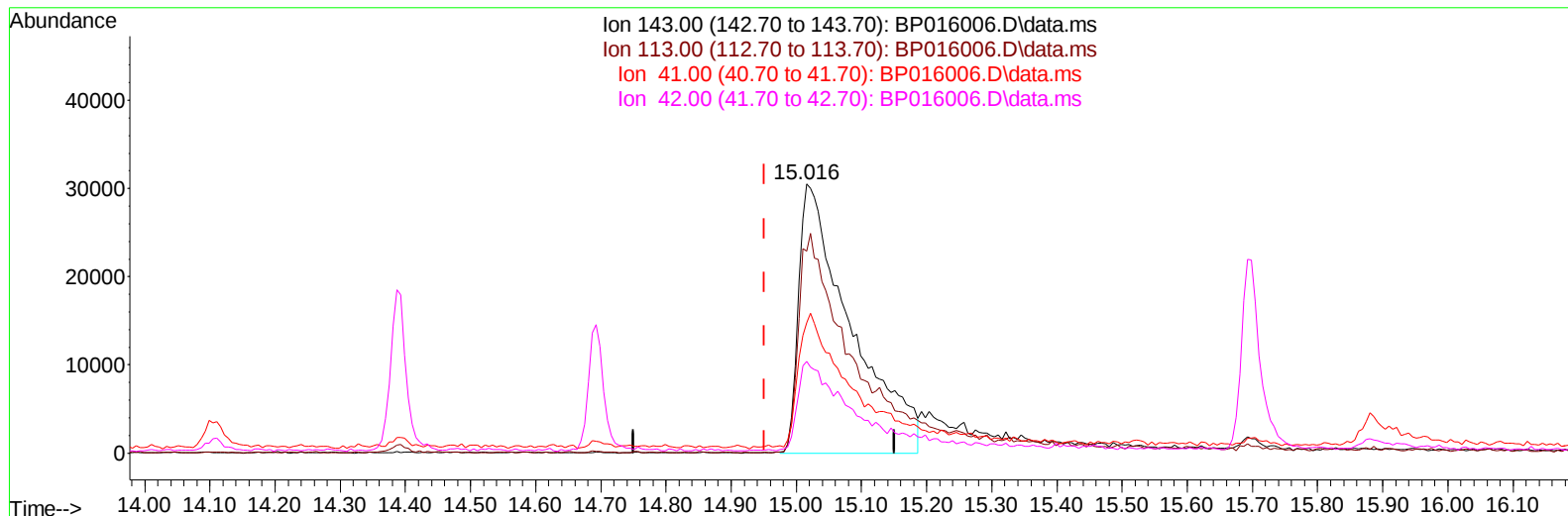
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TIC: BP016006.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.016min (+ 0.065) 27.51 ng/ul m

response 169032

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	75.18#
41.00	45.70	48.40
42.00	31.20	34.13

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
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 Operator : MA/JU
 Sample : PB153469BL
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK469

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/05/2023
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Quant Time: Jul 05 11:40:55 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	230870	20.000	ng/ul	0.00
20) Naphthalene-d8	10.869	136	1032443	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.692	164	602466	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.457	188	1092796	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	750397	20.000	ng/ul	# 0.00
88) Perylene-d12	24.098	264	840372	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	50708	7.902	ng/uL	0.00
4) Pyridine-d5	3.811	84	701360	43.359	ng/ul	0.00
7) Phenol-d5	7.228	99	894867	45.302	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	578504	47.337	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	721342	48.375	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	708792	45.421	ng/ul	0.00
21) Nitrobenzene-d5	9.228	128	332617	42.155	ng/ul	0.00
24) 2-Nitrophenol-d4	9.963	143	360351	39.130	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	591826	36.064	ng/ul	0.02
31) 4-Chloroaniline-d4	11.040	131	826033	39.184	ng/ul	0.01
46) Dimethylphthalate-d6	14.104	166	1756222	37.715	ng/ul	0.00
49) Acenaphthylene-d8	14.387	160	1984244	37.906	ng/ul	0.00
54) 4-Nitrophenol-d4	15.016	143	169032m	27.507	ng/ul	0.06
60) Fluorene-d10	15.698	176	1398863	36.483	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.886	200	107724	16.029	ng/ul	0.03
73) Anthracene-d10	17.563	188	1912710	38.318	ng/ul	0.00
81) Pyrene-d10	19.792	212	1865976	38.082	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264	1704952	40.219	ng/ul	0.00

Target Compounds Qvalue

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

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