

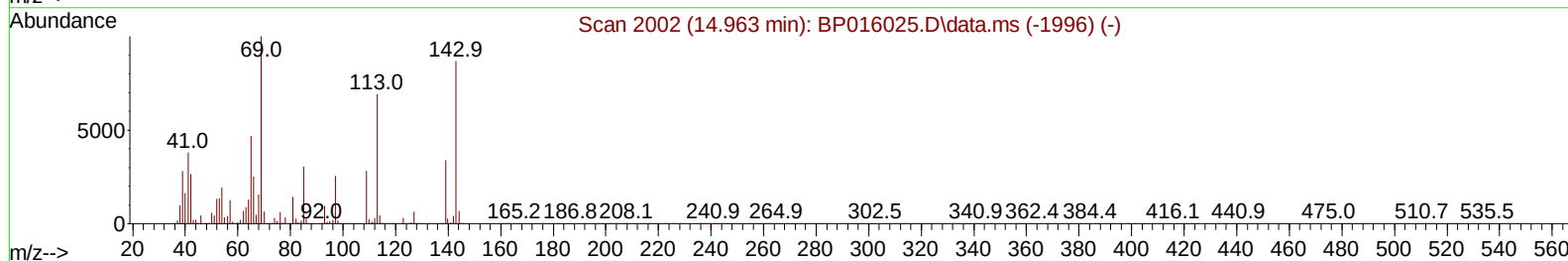
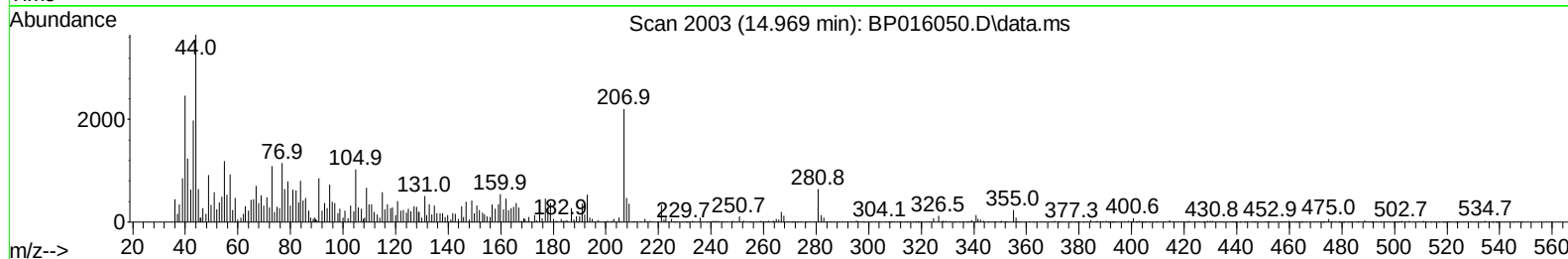
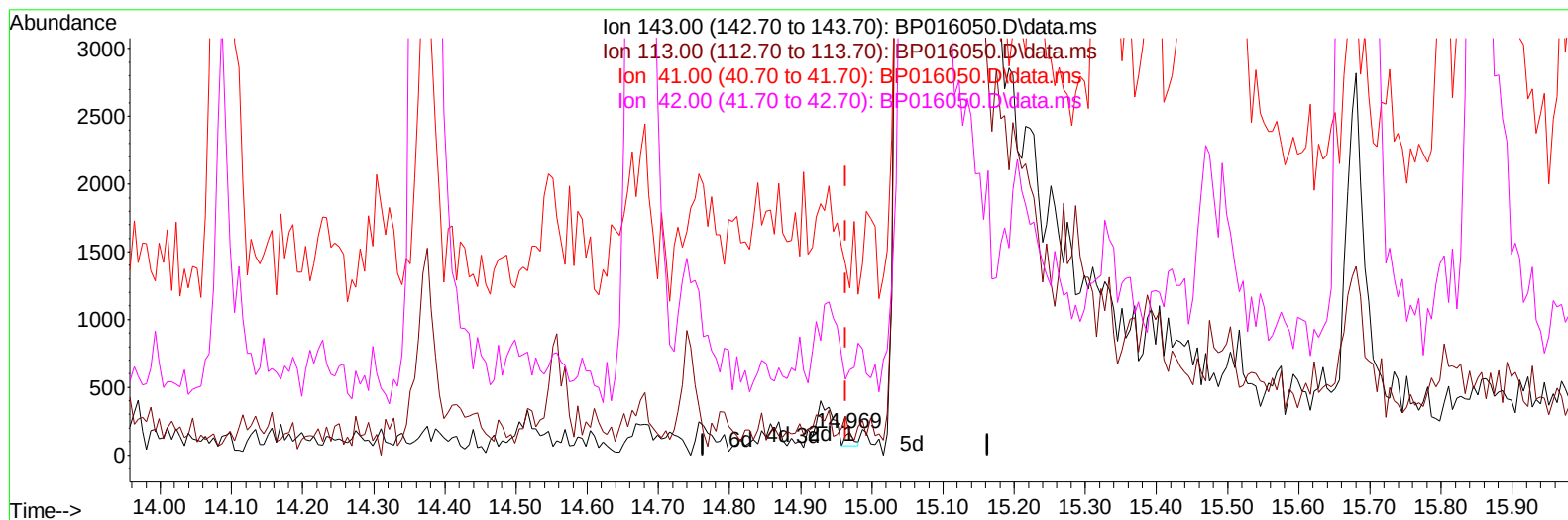
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070623\
 Data File : BP016050.D
 Acq On : 07 Jul 2023 01:28
 Operator : MA/JU
 Sample : 03284-08
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGZ5

Manual IntegrationsAPPROVED

Quant Time: Jul 07 02:25:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 06 23:18:26 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/07/2023
 Supervised By :mohammad ahmed 07/07/2023



TIC: BP016050.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.006) 0.01 ng/ul

response 82

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	90.36
41.00	45.70	742.77#
42.00	31.20	378.92#

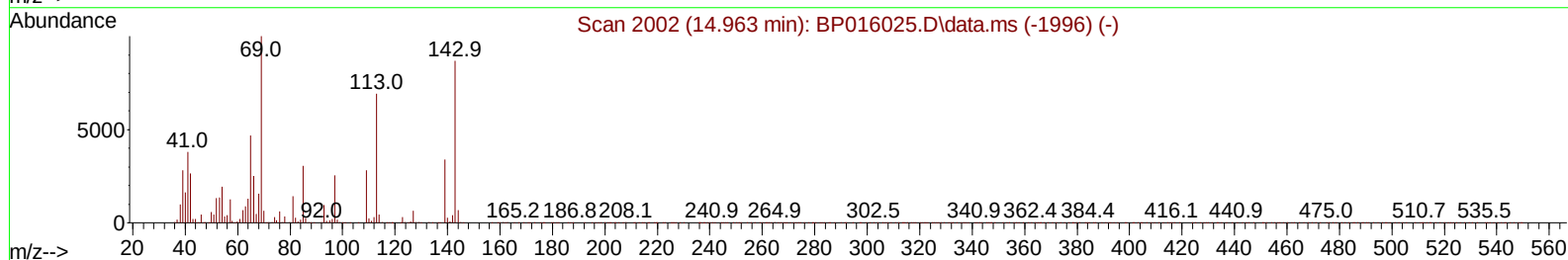
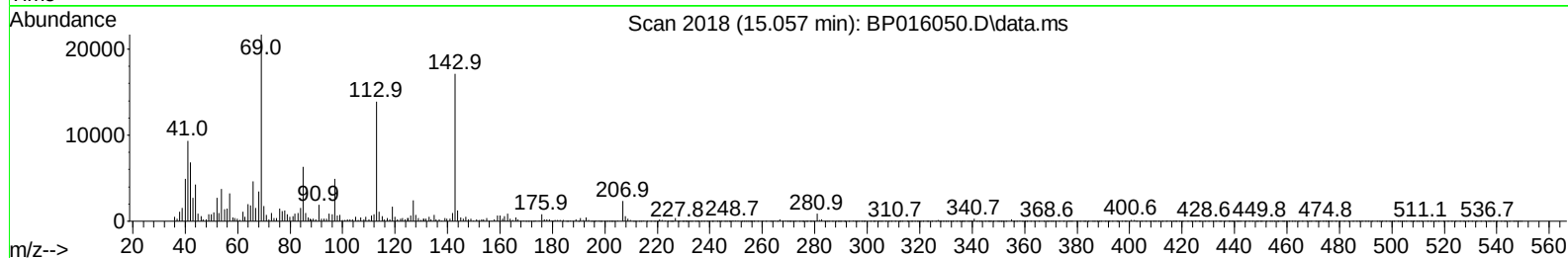
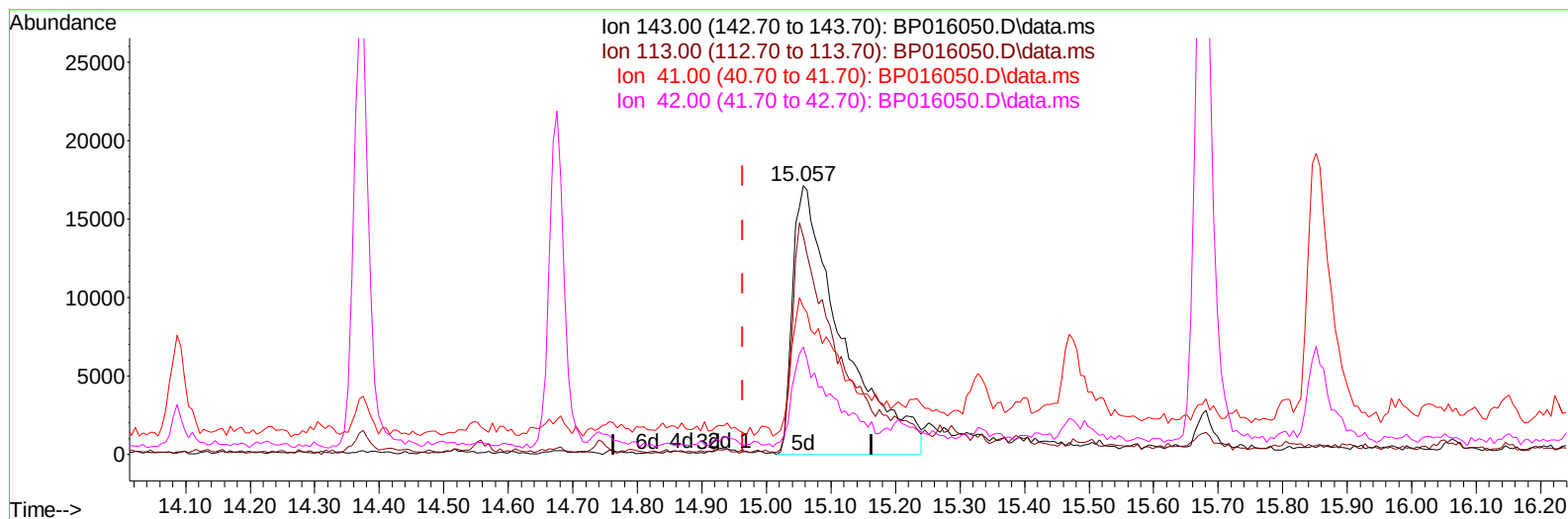
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Misc :
ALS Vial : 27 Sample Multiplier: 1

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Manual IntegrationsAPPROVED

Quant Time: Jul 07 02:26:33 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
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TIC: BP016050.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.057min (+ 0.094) 9.15 ng/ul m

response 90189

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	81.26
41.00	45.70	54.79
42.00	31.20	40.00#

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 Operator : MA/JU
 Sample : 03284-08
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGZ5

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/07/2023
 Supervised By :mohammad ahmed 07/07/2023

Quant Time: Jul 07 02:27:07 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	335964	20.000	ng/ul	0.00
20) Naphthalene-d8	10.852	136	1477746	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.675	164	966368	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	2122968	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1837867	20.000	ng/ul	0.00
88) Perylene-d12	24.074	264	1729099	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	45121	4.832	ng/uL	0.00
4) Pyridine-d5	3.817	84	142121	6.038	ng/ul	0.01
7) Phenol-d5	7.234	99	246513	8.576	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.352	67	660844	37.159	ng/ul	0.00
11) 2-Chlorophenol-d4	7.564	132	643042	29.634	ng/ul	0.00
15) 4-Methylphenol-d8	8.769	113	502422	22.125	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	407477	36.081	ng/ul	0.00
24) 2-Nitrophenol-d4	9.952	143	438195	33.245	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.516	165	765487	32.590	ng/ul	0.01
31) 4-Chloroaniline-d4	11.034	131	912382	30.238	ng/ul	0.01
46) Dimethylphthalate-d6	14.087	166	3067483	41.068	ng/ul	0.00
49) Acenaphthylene-d8	14.375	160	3168948	37.741	ng/ul	0.00
54) 4-Nitrophenol-d4	15.057	143	90189m	9.150	ng/ul	0.09
60) Fluorene-d10	15.675	176	2463571	40.057	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	470072	36.004	ng/ul	0.00
73) Anthracene-d10	17.545	188	3824510	39.439	ng/ul	0.00
81) Pyrene-d10	19.769	212	4790341	39.917	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.904	264	3544724	40.640	ng/ul	0.00

Target Compounds Qvalue

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

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