

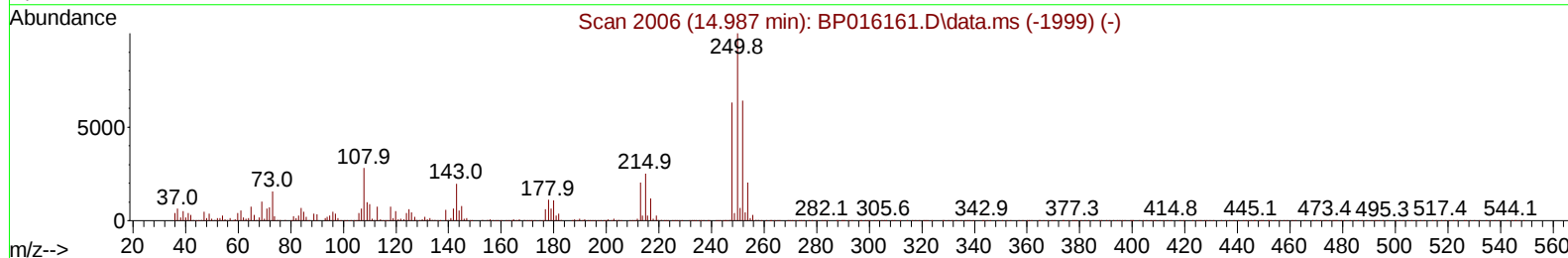
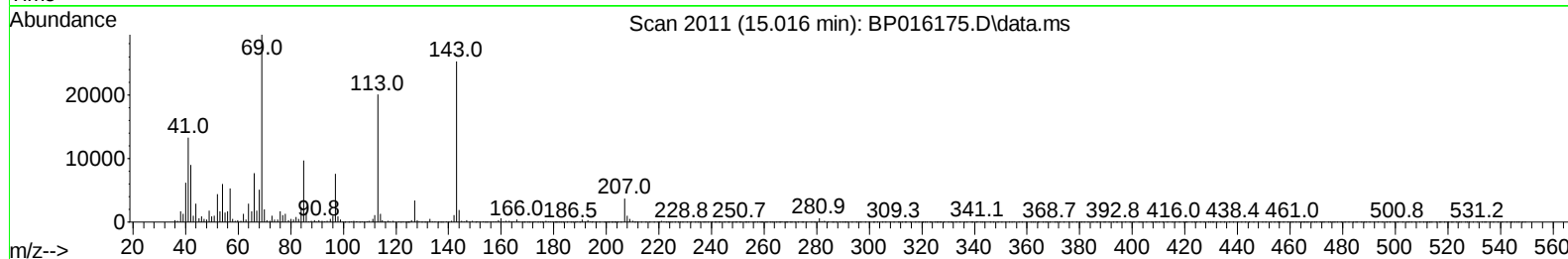
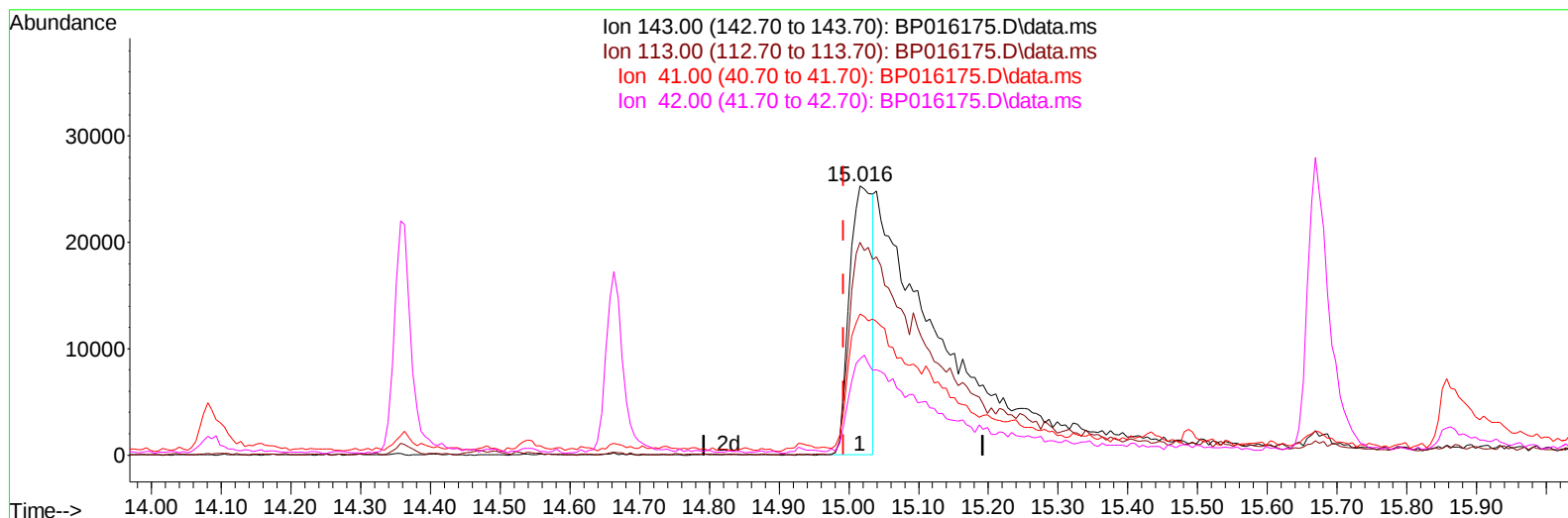
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071023\
Data File : BP016175.D
Acq On : 11 Jul 2023 02:02
Operator : MA/JU
Sample : 03404-16
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YBW52

Manual IntegrationsAPPROVED

Quant Time: Jul 11 03:09:41 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 10 23:19:51 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/11/2023
Supervised By :mohammad ahmed 07/11/2023



TIC: BP016175.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.016min (+ 0.024) 7.85 ng/u1

response 57613

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	79.21
41.00	45.70	52.51
42.00	31.20	35.65

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Quant Time: Jul 11 03:26:07 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	238710	20.000	ng/ul	0.00
20) Naphthalene-d8	10.840	136	1115343	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.663	164	719564	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1576504	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.545	240	1350056	20.000	ng/ul	# 0.00
88) Perylene-d12	24.068	264	1268807	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	34336	5.175	ng/uL	0.00
4) Pyridine-d5	3.793	84	337072	20.154	ng/ul	0.00
7) Phenol-d5	7.211	99	651527	31.899	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	472319	37.379	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	518719	33.644	ng/ul	0.00
15) 4-Methylphenol-d8	8.763	113	571765	35.436	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	274445	32.197	ng/ul	0.00
24) 2-Nitrophenol-d4	9.952	143	311154	31.277	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.540	165	440402	24.842	ng/ul	0.03
31) 4-Chloroaniline-d4	11.057	131	490390	21.534	ng/ul	0.04
46) Dimethylphthalate-d6	14.081	166	2108267	37.907	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	2317592	37.069	ng/ul	0.00
54) 4-Nitrophenol-d4	15.016	143	188688m	25.709	ng/ul	0.02
60) Fluorene-d10	15.669	176	1799964	39.305	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	221944	22.892	ng/ul	0.00
73) Anthracene-d10	17.534	188	2836694	39.392	ng/ul	0.00
81) Pyrene-d10	19.763	212	3345634	37.951	ng/ul	-0.02
92) Benzo(a)pyrene-d12	23.898	264	2474125	38.656	ng/ul	-0.02

Target Compounds Qvalue

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

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