

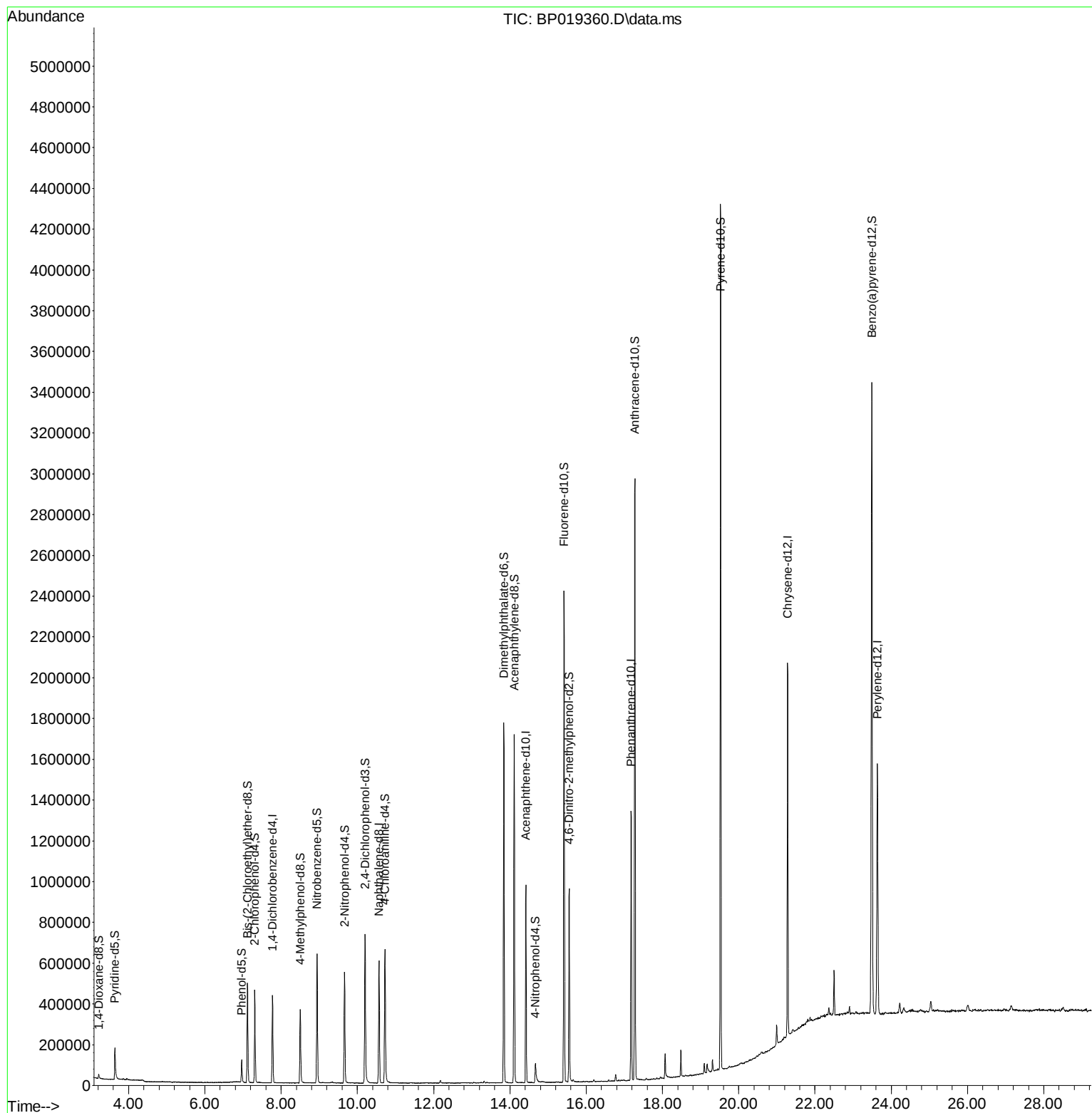
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019360.D
Acq On : 13 Jan 2024 08:51
Operator : MA/JU
Sample : P1130-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
Supervised By :mohammad ahmed 01/16/2024

Quant Time: Jan 14 20:33:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jan 09 03:05:24 2024
Response via : Initial Calibration



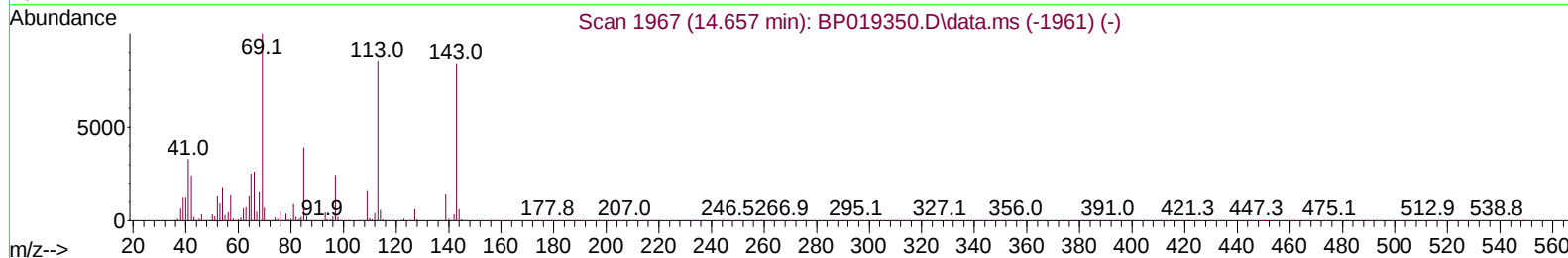
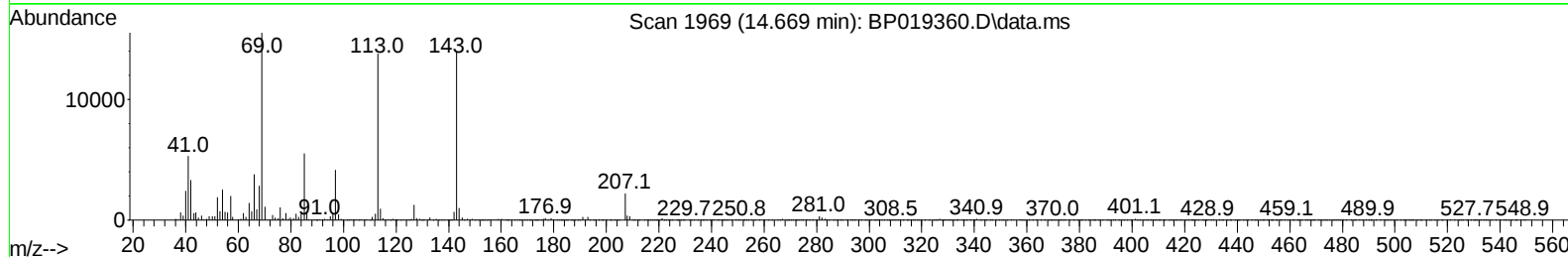
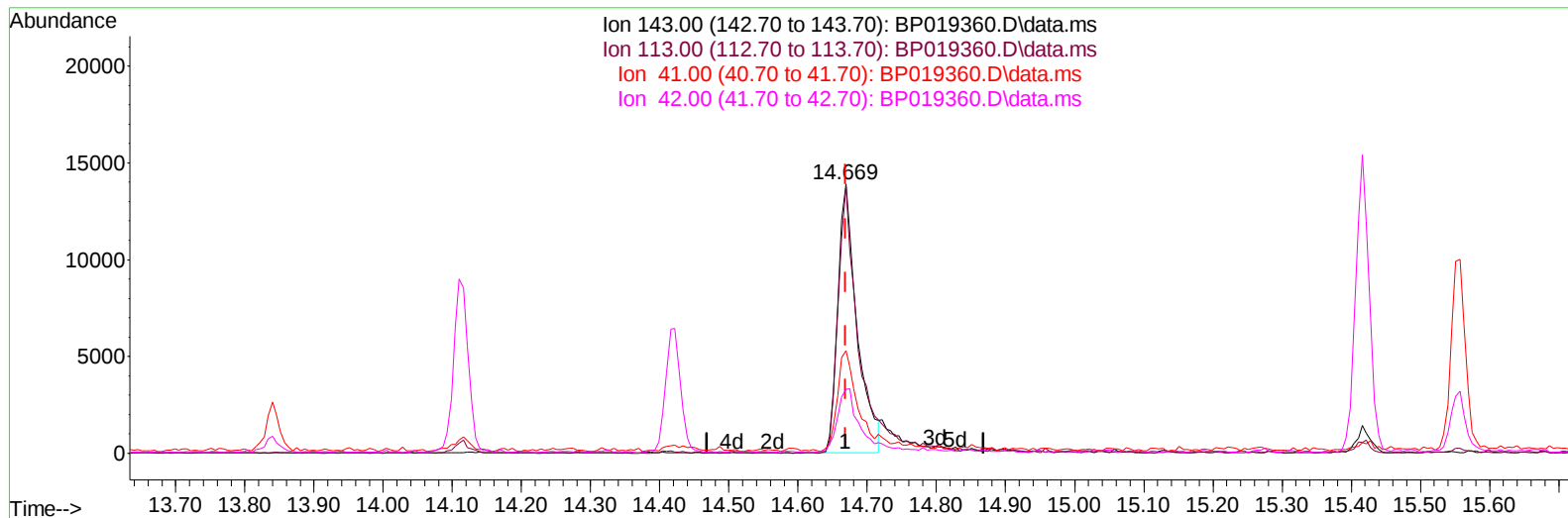
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TIC: BP019360.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 6.29 ng/u1

response 26967

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	99.14
41.00	39.60	38.00
42.00	25.70	23.78

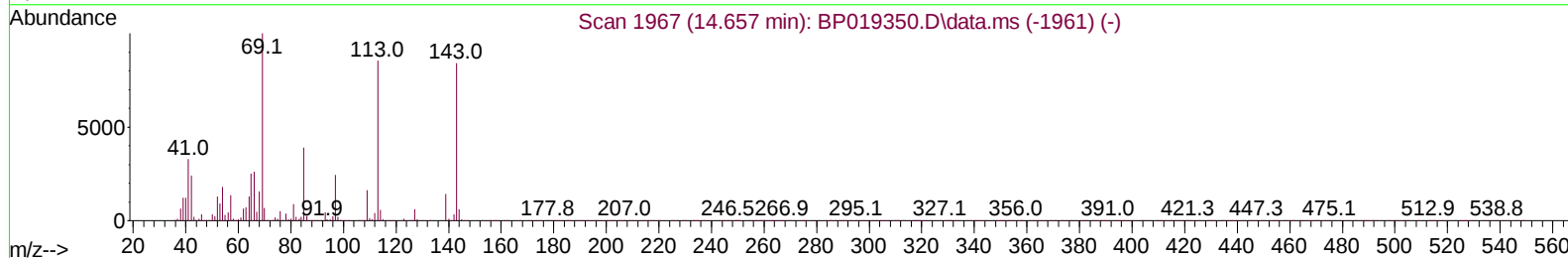
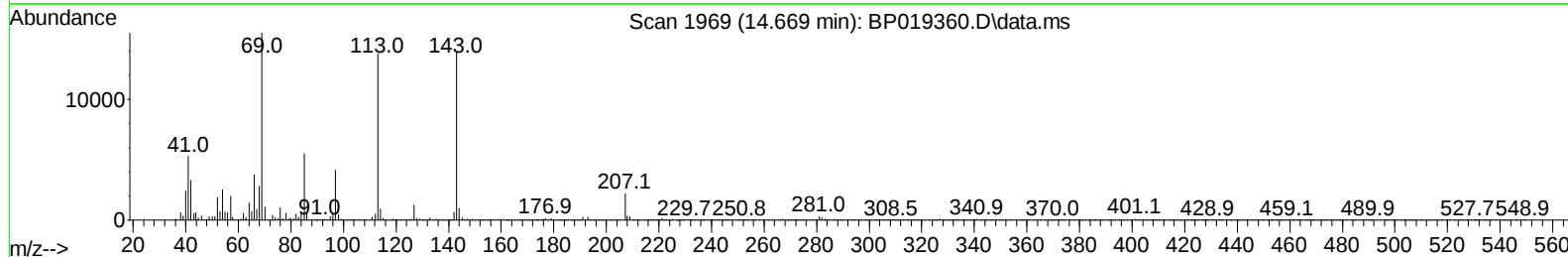
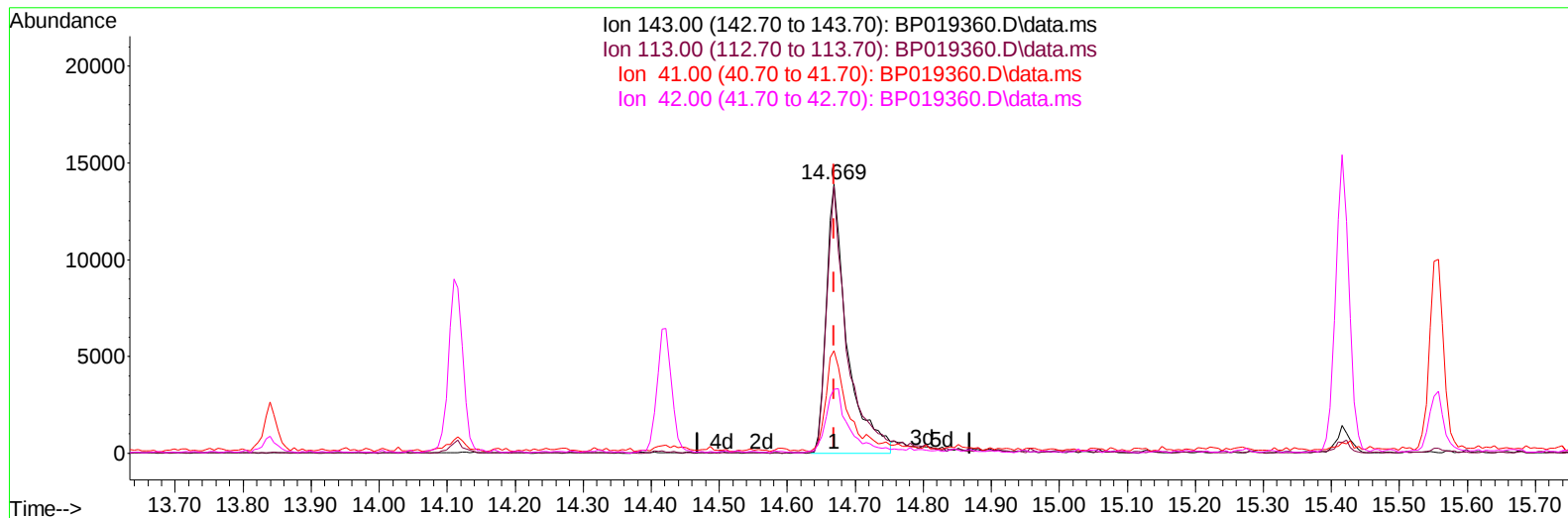
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Instrument :
 BNA_P
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TIC: BP019360.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 6.84 ng/ul m

response 29332

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	99.14
41.00	39.60	38.00
42.00	25.70	23.78

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
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 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AY5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	109802	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.569	136	445517	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.422	164	315899	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.175	188	815855	20.000	ng/ul	-0.01
79) Chrysene-d12	21.280	240	890254	20.000	ng/ul	0.00
88) Perylene-d12	23.639	264	964237	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	12039	4.146	ng/uL	0.00
4) Pyridine-d5	3.640	84	93898	11.183	ng/ul	0.00
7) Phenol-d5	6.963	99	63904	6.089	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	216158	34.435	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.311	132	182213	25.276	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	132089	15.865	ng/ul	-0.01
21) Nitrobenzene-d5	8.940	128	131495	35.083	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.663	143	150397	34.534	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	272241	30.942	ng/ul	0.00
31) 4-Chloroaniline-d4	10.722	131	314532	28.301	ng/ul	0.00
46) Dimethylphthalate-d6	13.840	166	1119961	40.129	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	1074082	35.902	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	29332m	6.844	ng/ul	0.00
60) Fluorene-d10	15.416	176	928035	40.615	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	229051	38.027	ng/ul	0.00
73) Anthracene-d10	17.281	188	1730869	42.538	ng/ul	0.00
81) Pyrene-d10	19.522	212	2334375	40.213	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.492	264	2357054	44.956	ng/ul	0.00

Target Compounds Qvalue

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