

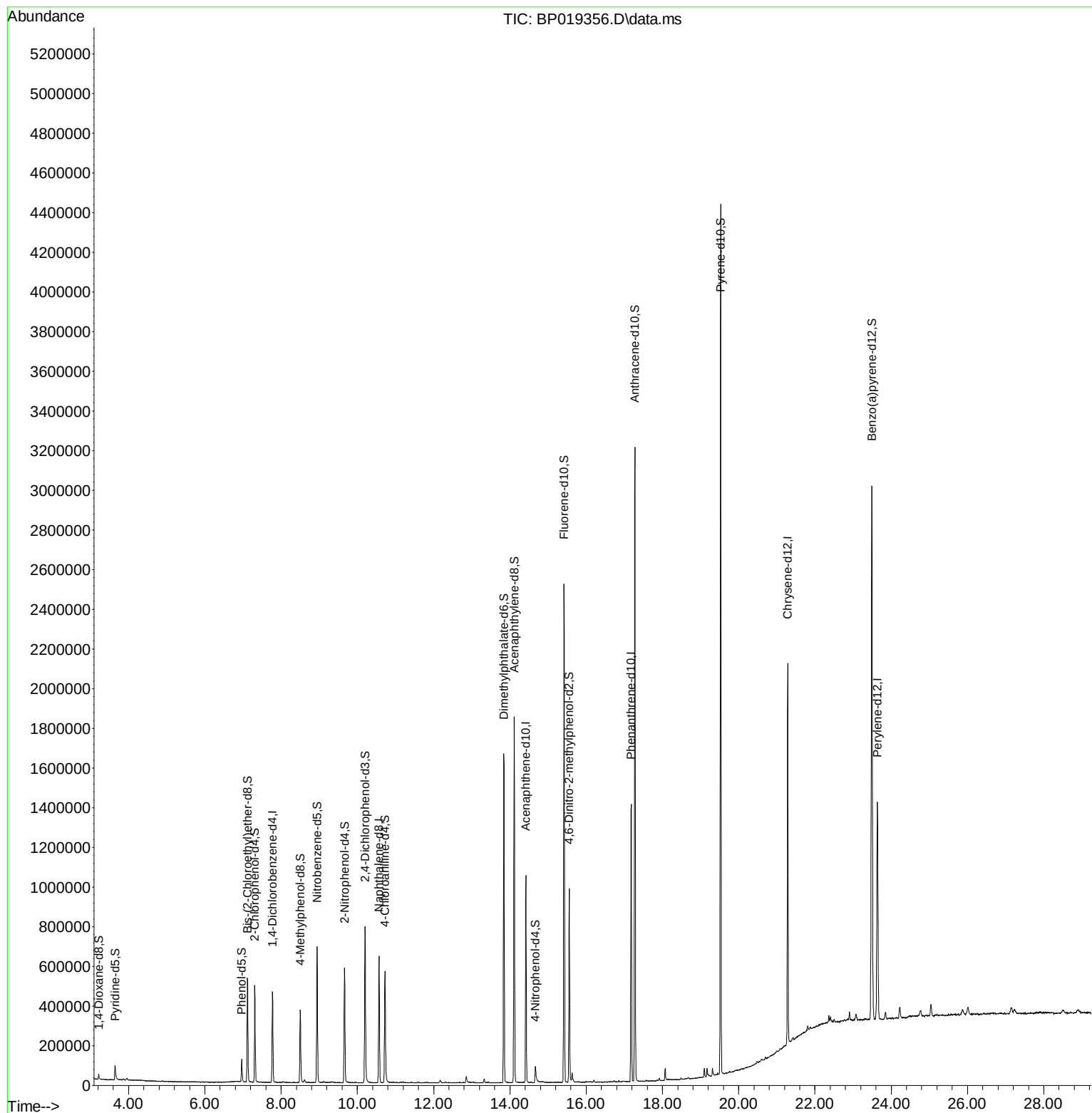
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019356.D
Acq On : 13 Jan 2024 06:35
Operator : MA/JU
Sample : P1091-19
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AZ9

Manual IntegrationsAPPROVED

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

Quant Time: Jan 14 20:25:02 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



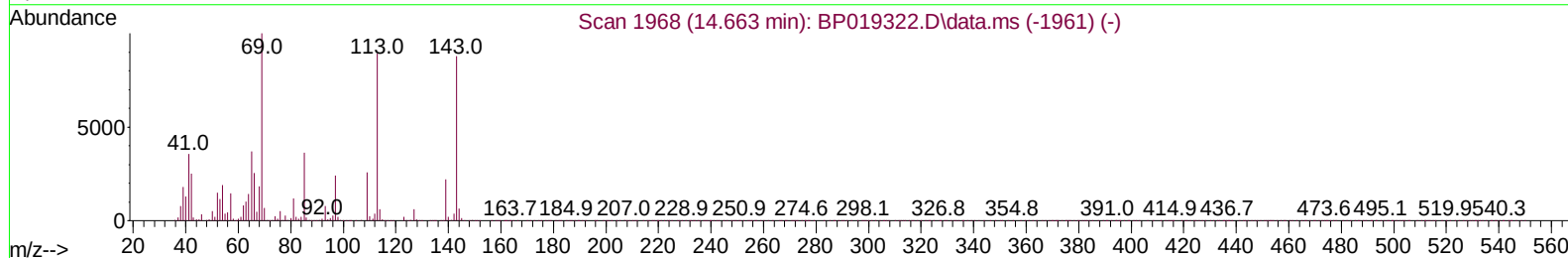
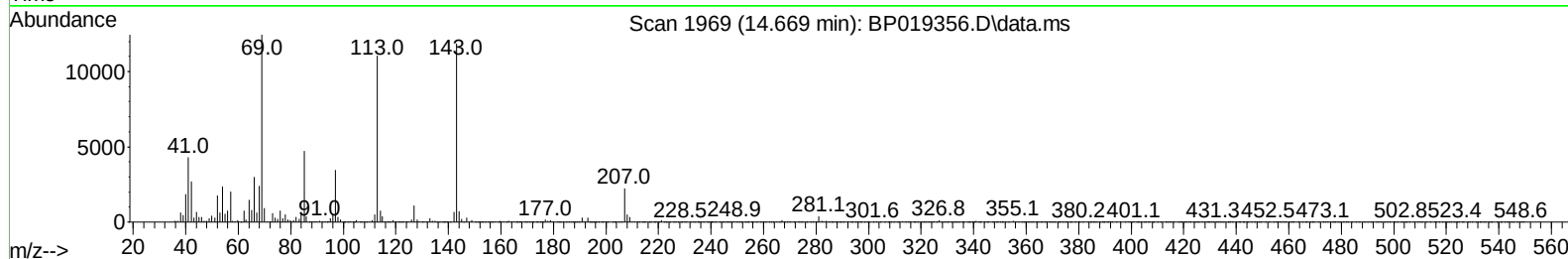
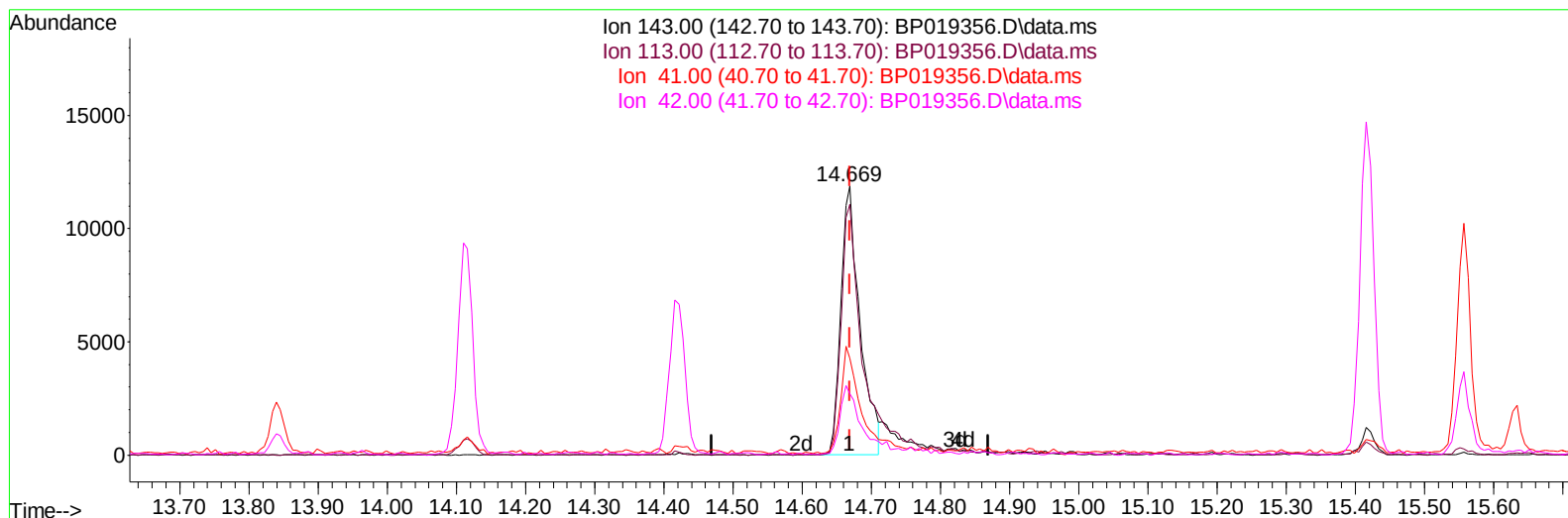
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Quant Time: Jan 14 20:22:18 2024
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TIC: BP019356.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 4.99 ng/u1

response 22887

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	93.17
41.00	39.60	36.08
42.00	25.70	22.78

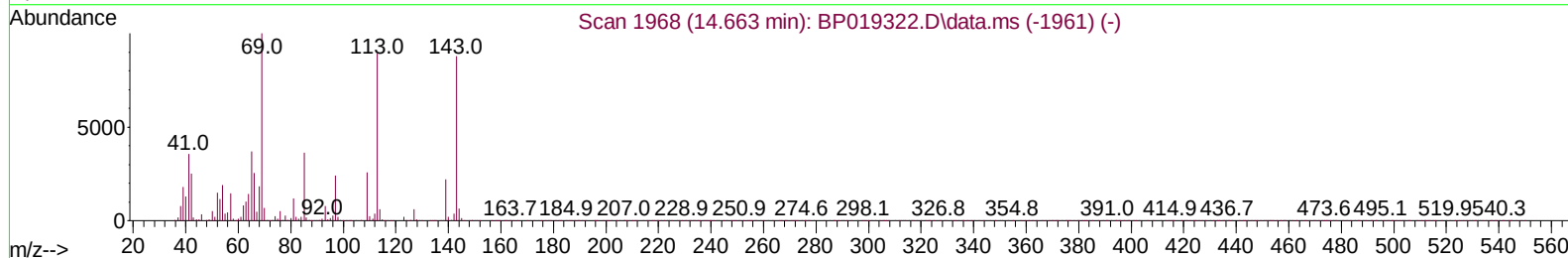
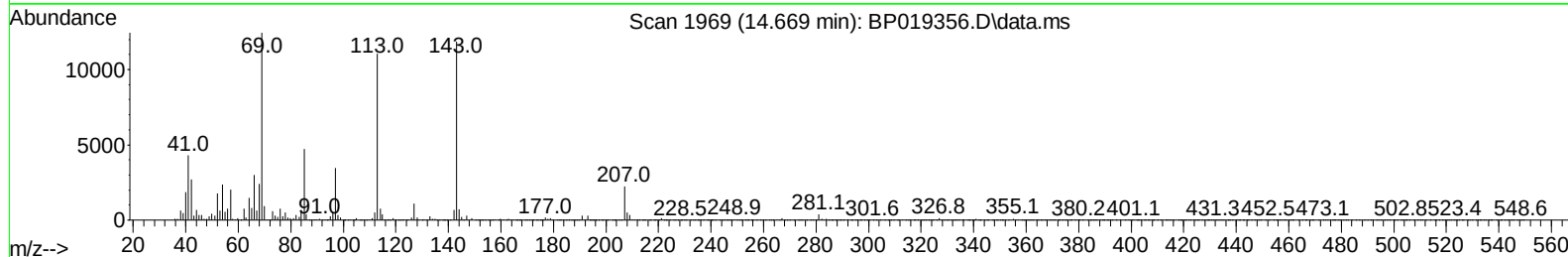
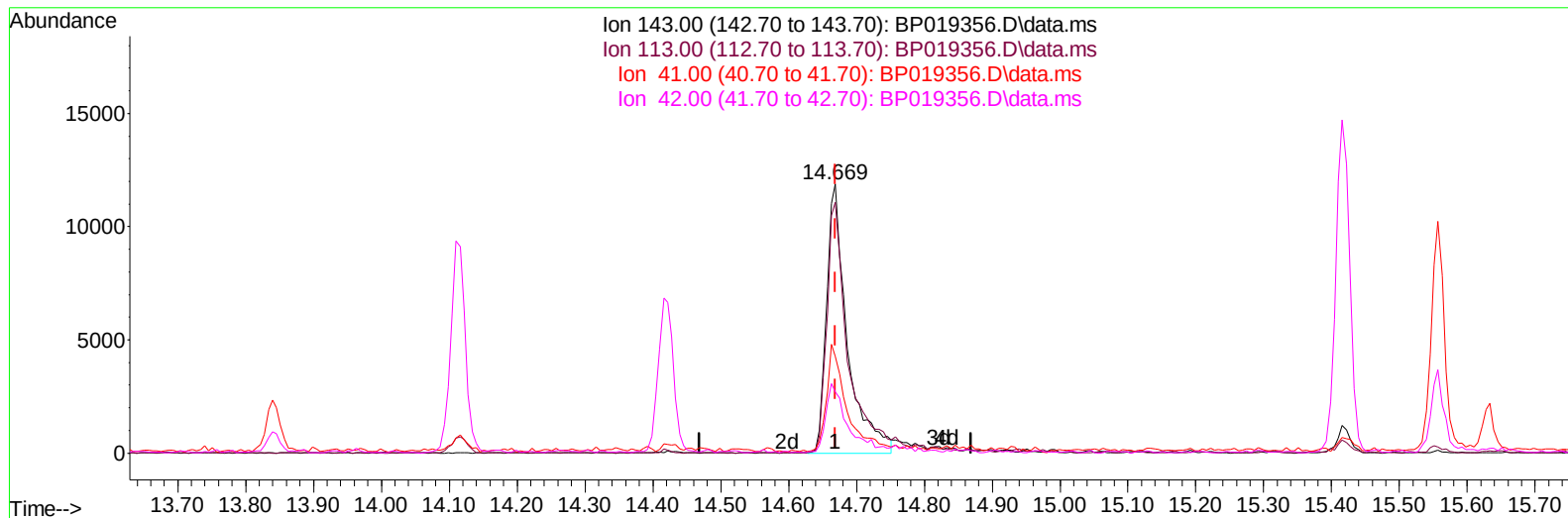
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TIC: BP019356.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 5.51 ng/ul m

response 25265

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	93.17
41.00	39.60	36.08
42.00	25.70	22.78

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
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 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AZ9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 09 03:05:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	118552	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.569	136	471332	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.422	164	337793	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.180	188	848273	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	879602	20.000	ng/ul	0.00
88) Perylene-d12	23.639	264	911207	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	12991	4.144	ng/uL	0.00
4) Pyridine-d5	3.646	84	46462	5.125	ng/ul	0.00
7) Phenol-d5	6.963	99	64036	5.651	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	235732	34.782	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.310	132	194341	24.968	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	129089	14.360	ng/ul	-0.01
21) Nitrobenzene-d5	8.946	128	141042	35.569	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	161999	35.161	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	283482	30.455	ng/ul	0.00
31) 4-Chloroaniline-d4	10.722	131	275856	23.461	ng/ul	0.00
46) Dimethylphthalate-d6	13.845	166	1065448	35.702	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	1121035	35.043	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	25265m	5.513	ng/ul	0.00
60) Fluorene-d10	15.416	176	946586	38.742	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	222783	35.573	ng/ul	0.00
73) Anthracene-d10	17.280	188	1767746	41.784	ng/ul	0.00
81) Pyrene-d10	19.527	212	2285365	39.845	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.492	264	2154288	43.480	ng/ul	0.00

Target Compounds Qvalue

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