Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010299.D
Acq On : 11 May 2022 23:20
Operator : CG/JU
Sample : N2707-12
Misc :
ALS Vial : 22 Sample Multiplier: 1

Quant Time: May 12 01:16:23 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 12 01:10:15 2022
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
EXLA3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022
Supervised By :Yogesh Patel 05/17/2022

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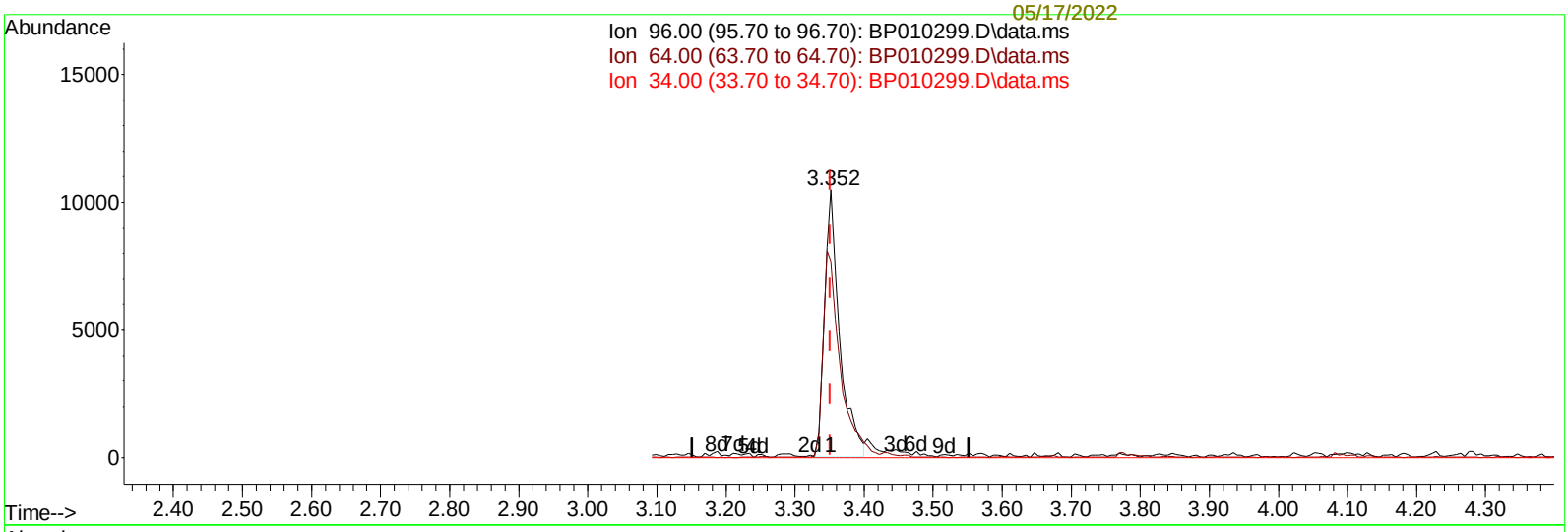
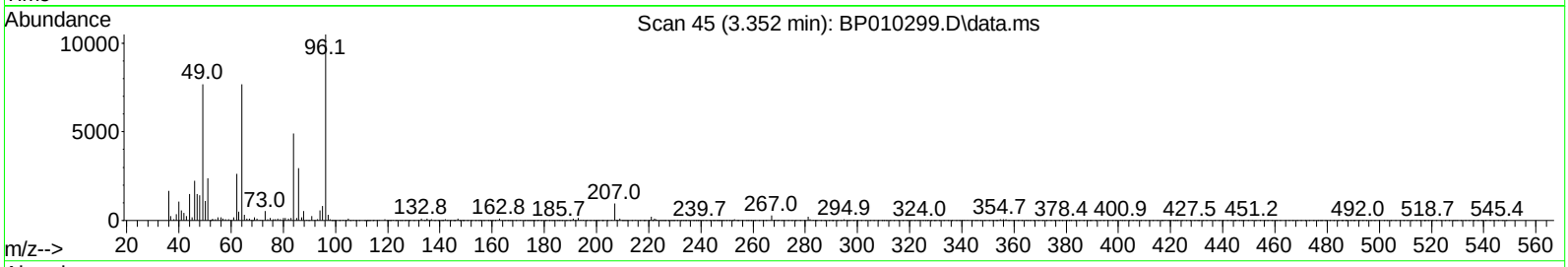
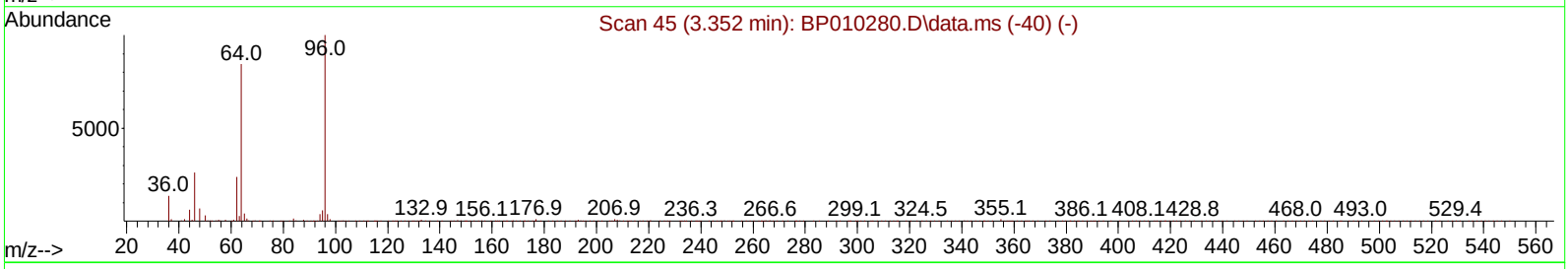
05/17/2022







TIC: BP010299.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.352min (+ 0.000) 4.13 ng/uL

response 15935

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.22#
34.00	0.00	0.00
0.00	0.00	0.00

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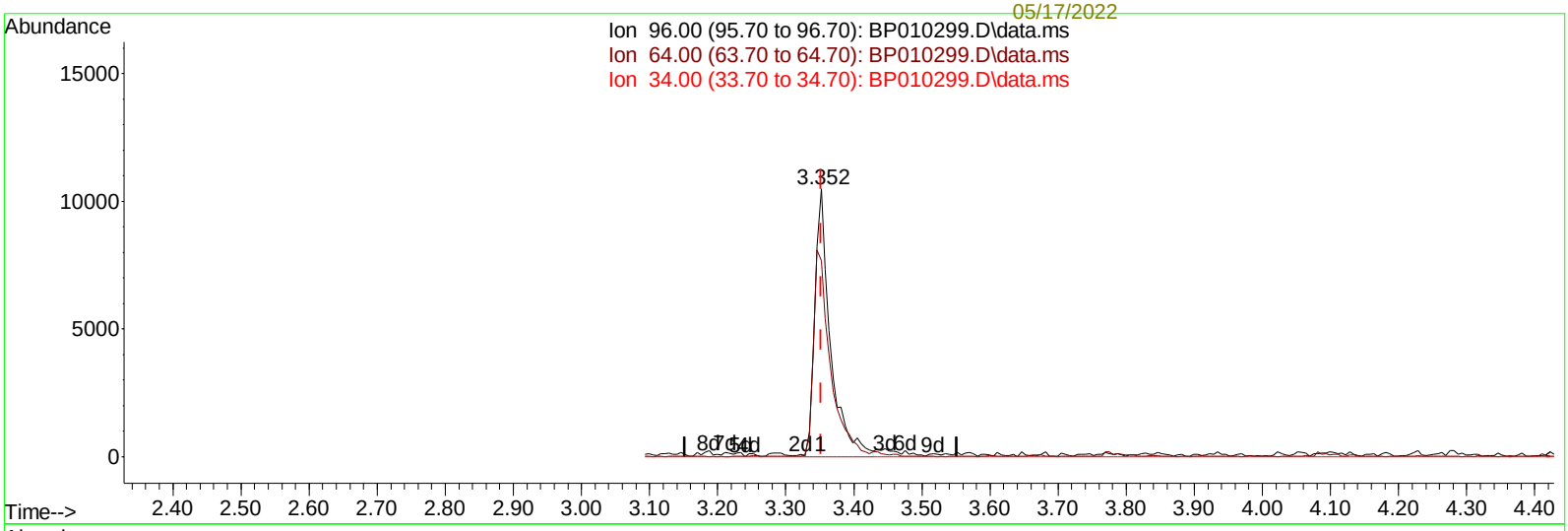

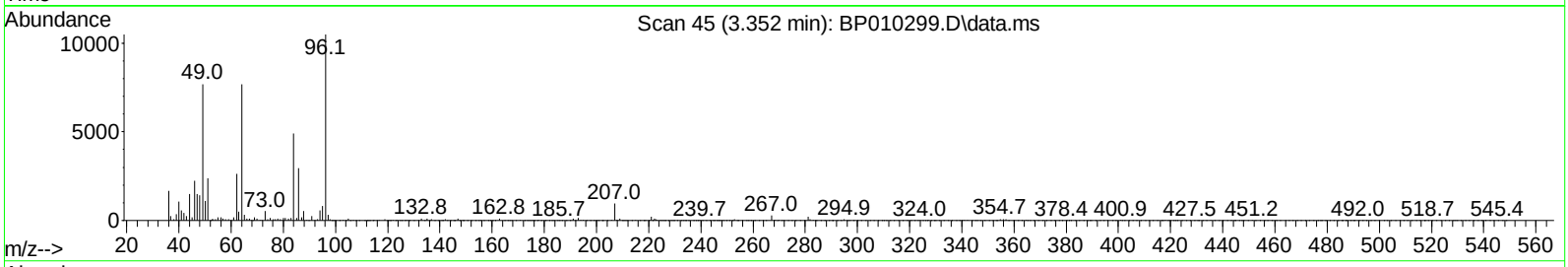

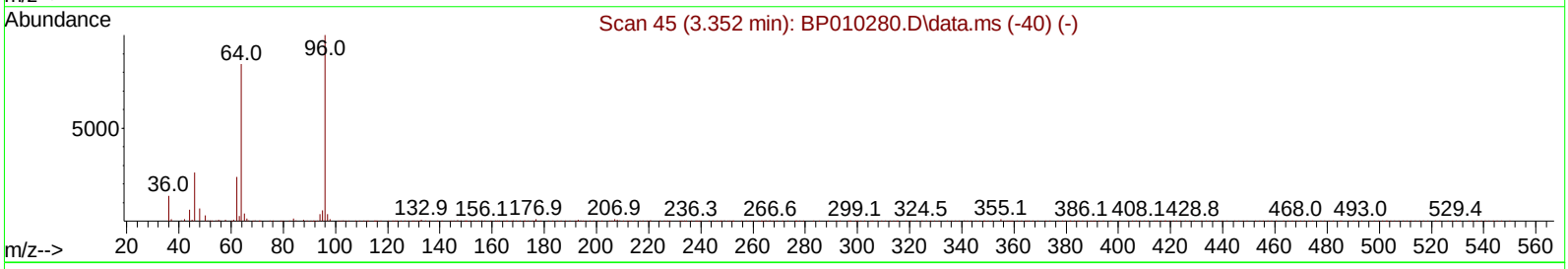

TIC: BP010299.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.352min (+ 0.000) 4.34 ng/uL m

response 16744

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.22#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
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ALS Vial : 22 Sample Multiplier: 1

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BNA_P
ClientSampleId :
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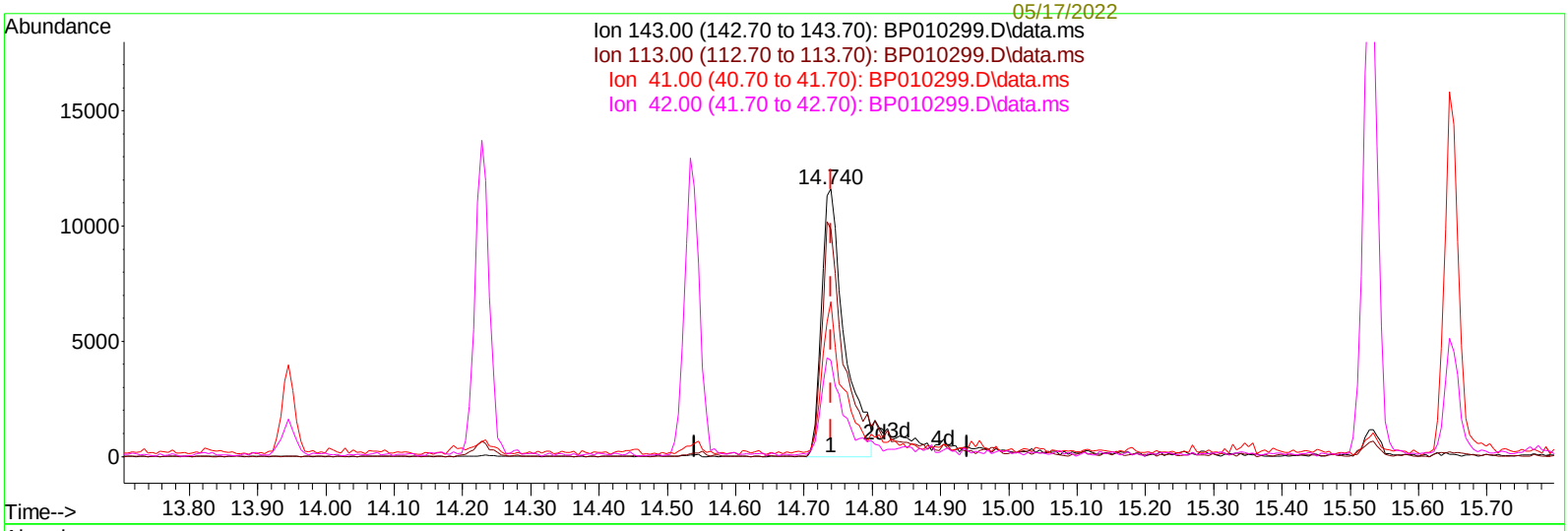

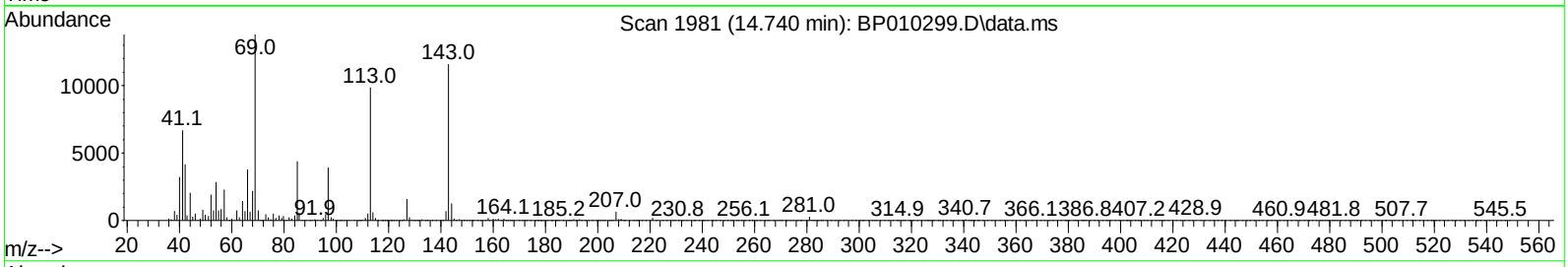

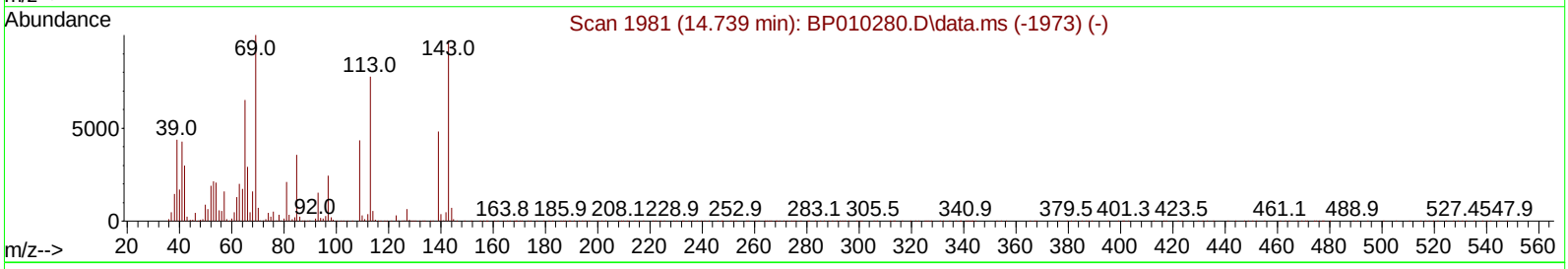

TIC: BP010299.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.000) 3.40 ng/ul

response 27165

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.09#
41.00	23.20	57.82#
42.00	18.00	35.95#

Quantitation Report (Qedit)

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Instrument :
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ClientSampleId :
EXLA3

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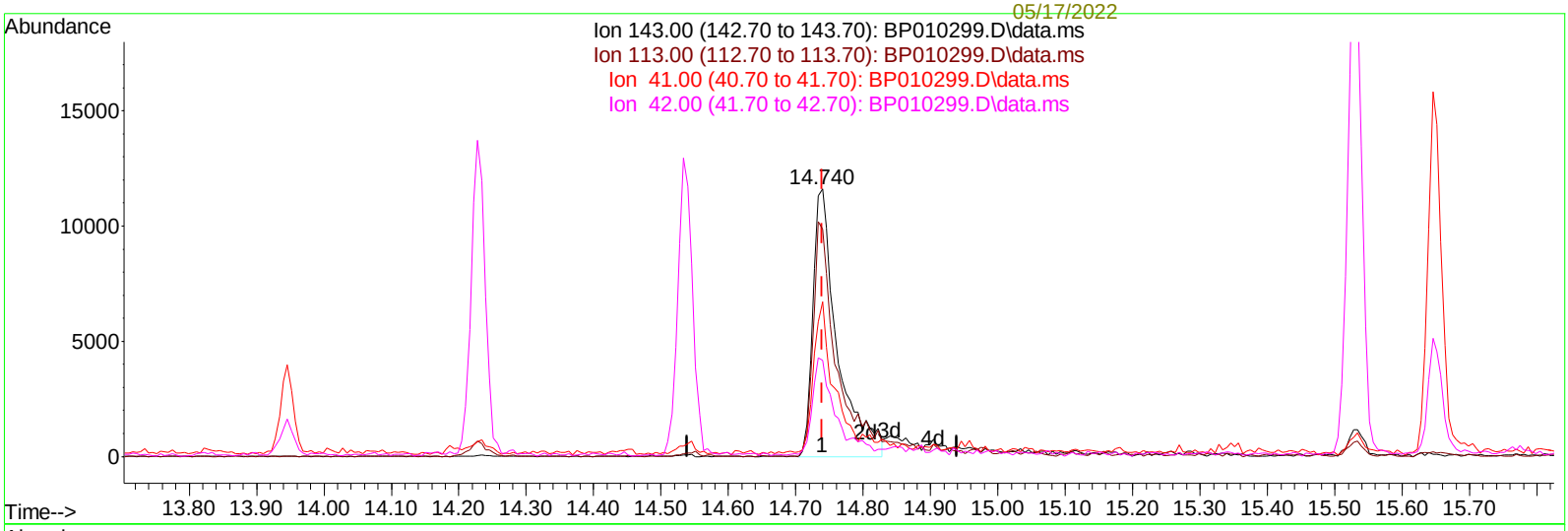

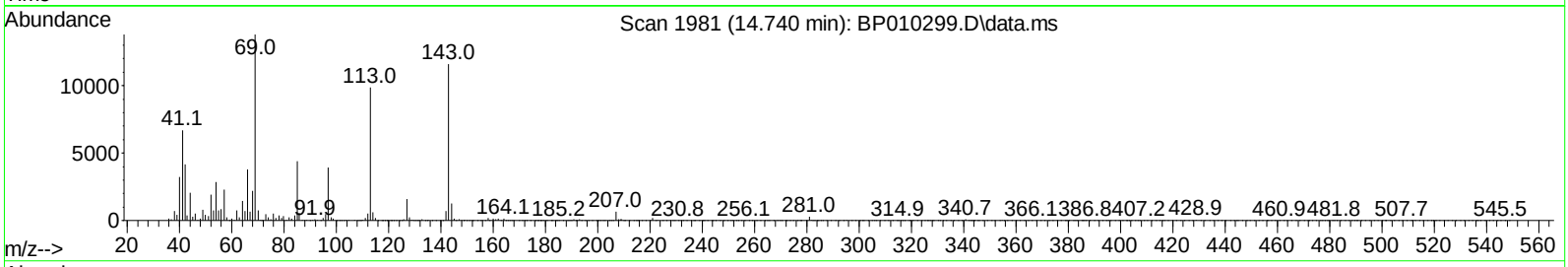

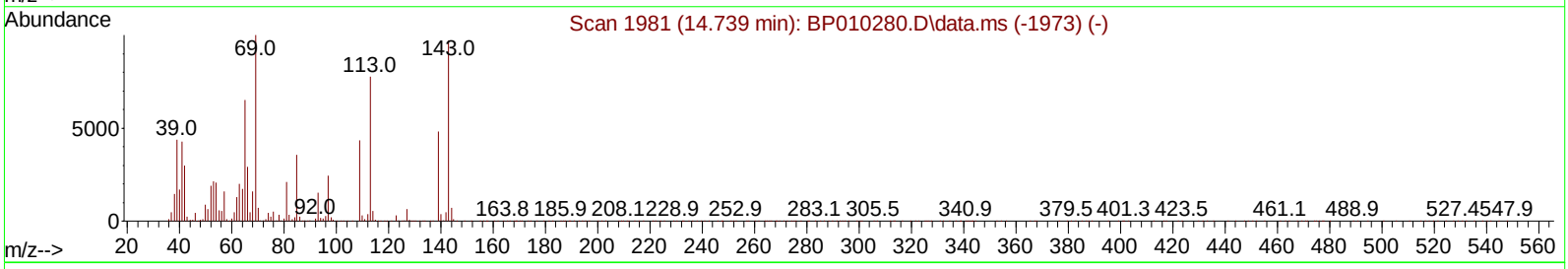

TIC: BP010299.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.000) 3.66 ng/ul m

response 29192

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.09#
41.00	23.20	57.82#
42.00	18.00	35.95#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	163781	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	735888	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.540	164	522443	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.287	188	1139046	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	1115194	20.000	ng/ul	# 0.00
88) Perylene-d12	23.804	264	1088650	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	16744m	4.343	ng/uL	0.00
4) Pyridine-d5	3.775	84	87848	8.188	ng/ul	0.00
7) Phenol-d5	7.069	99	72312	4.754	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	284445	30.227	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	243360	21.723	ng/ul	0.00
15) 4-Methylphenol-d8	8.611	113	158842	12.701	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	175898	29.811	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	186631	27.741	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	300981	24.580	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	424597	23.488	ng/ul	0.00
46) Dimethylphthalate-d6	13.946	166	1315115	31.597	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1433504	28.888	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	29192m	3.656	ng/ul	0.00
60) Fluorene-d10	15.528	176	1130091	32.166	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	217562	27.892	ng/ul	0.00
73) Anthracene-d10	17.387	188	1934175	34.675	ng/ul	0.00
81) Pyrene-d10	19.622	212	2310674	37.495	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	2095595	36.095	ng/ul	-0.01
Target Compounds						Qvalue

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

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