Quantitation Report (Qedit)

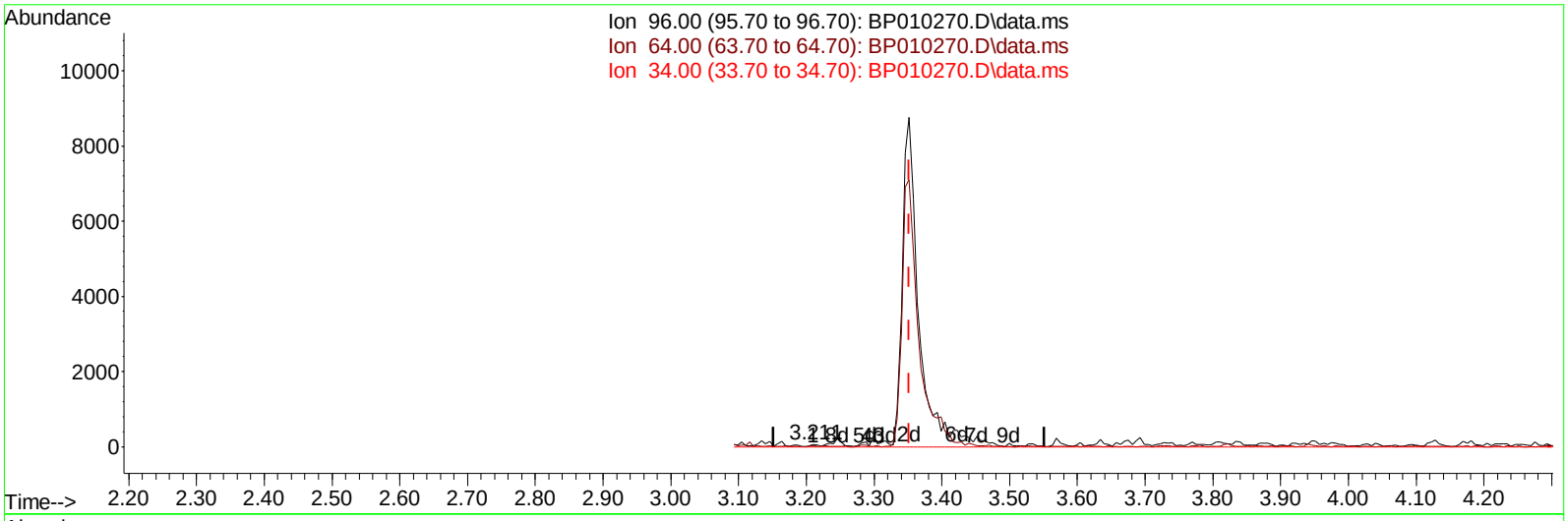
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051022\
Data File : BP010270.D
Acq On : 10 May 2022 17:59
Operator : CG/JU
Sample : N2678-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

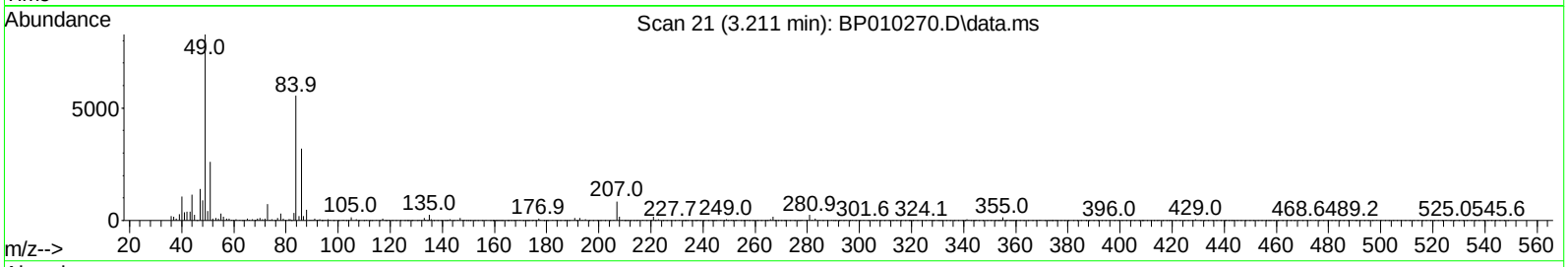
Quant Time: May 11 01:17:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 10 15:59:46 2022
Response via : Initial Calibration

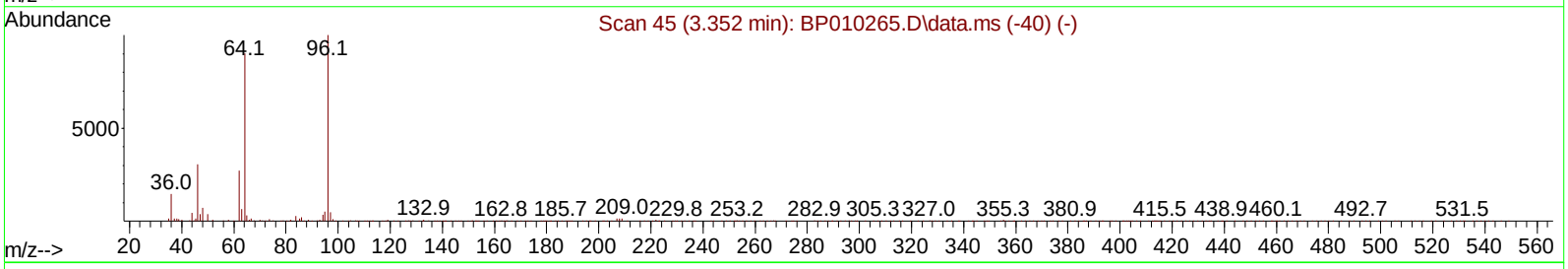
Instrument :
BNA_P
ClientSampleId :
BH0C1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/11/2022
Supervised By :mohammad ahmed 05/12/2022







TIC: BP010270.D\data.ms

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

3.211min (-0.141) 0.02 ng/uL

response 45

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	48.33
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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Misc :
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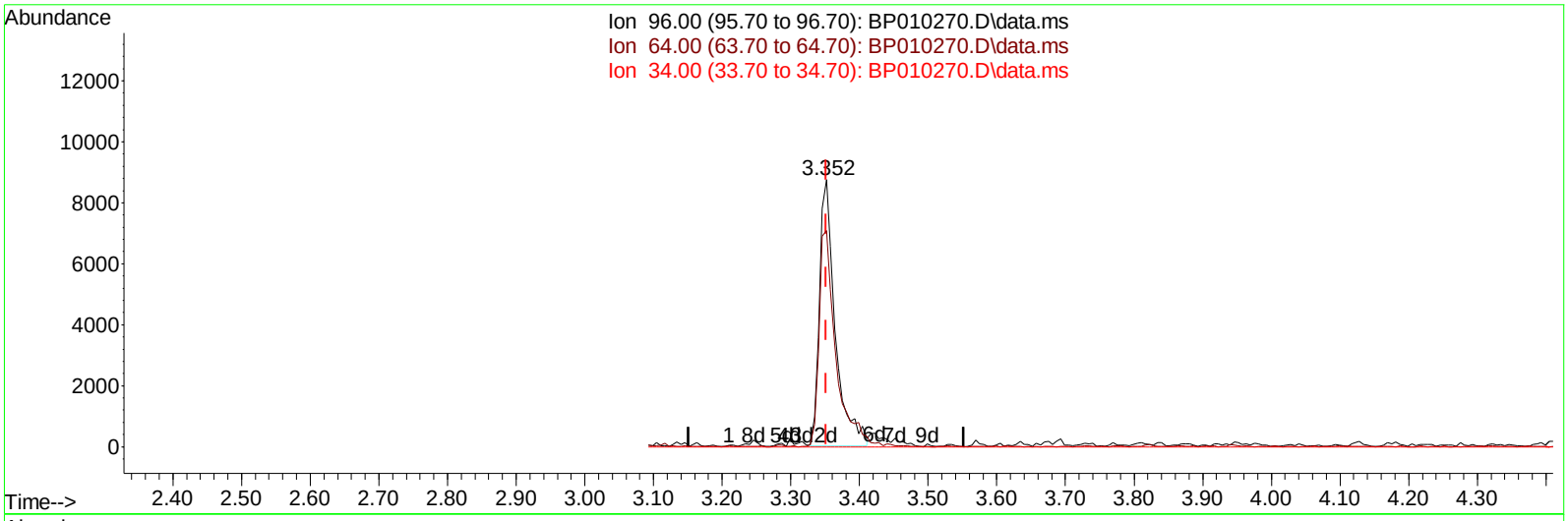

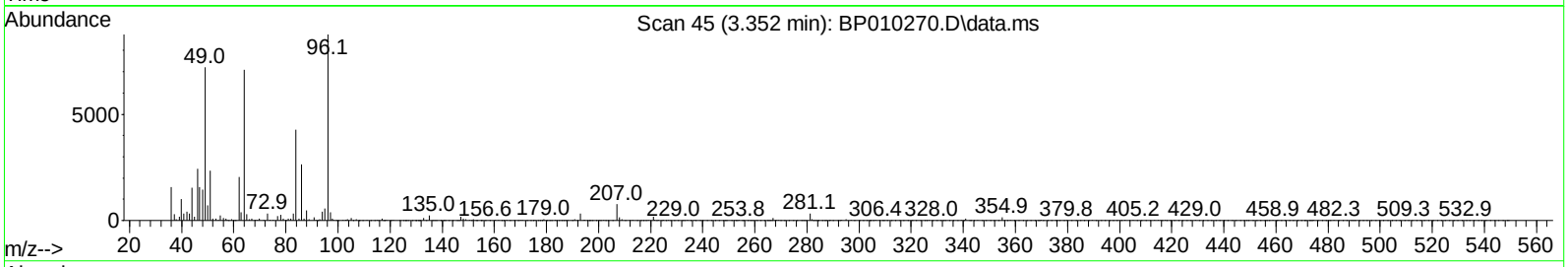

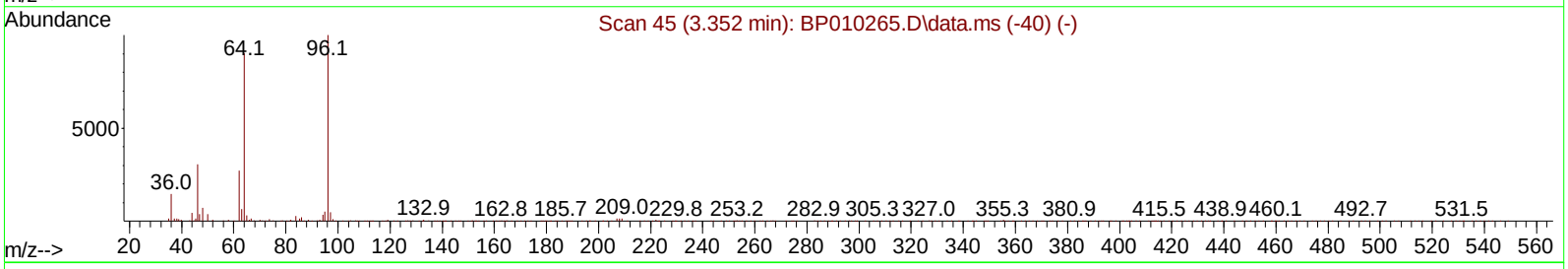

TIC: BP010270.D\data.ms

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

3.352min (-0.000) 5.73 ng/uL m

response 13885

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	81.10#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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Misc :
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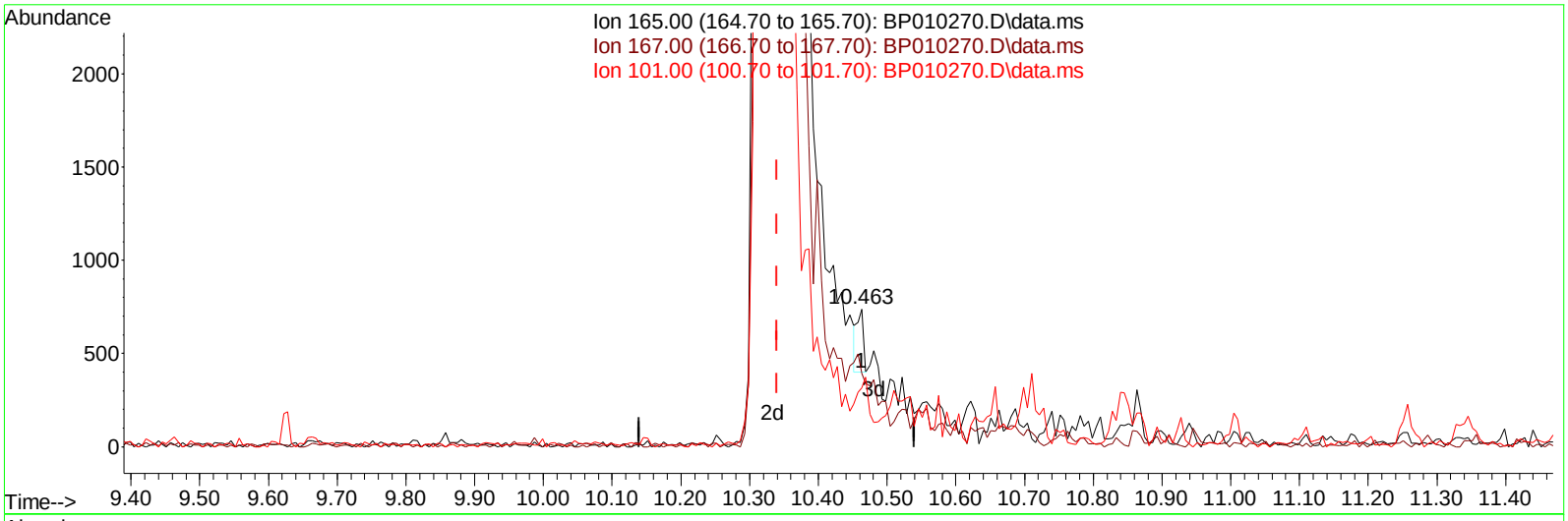

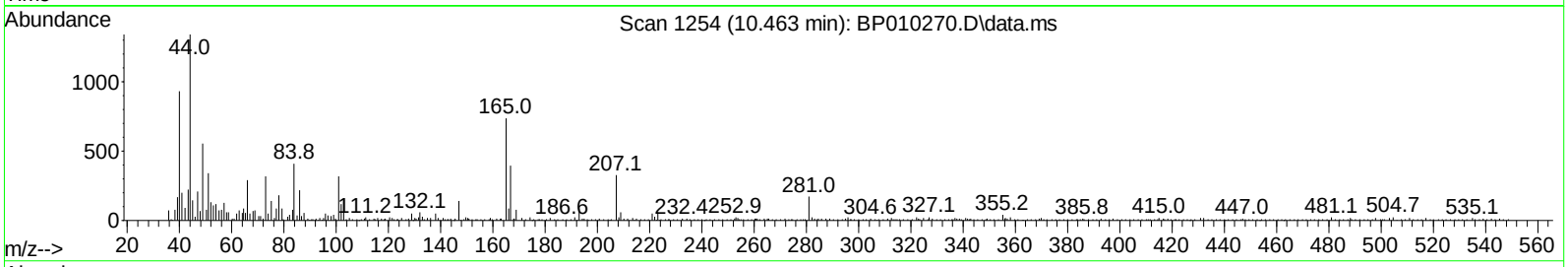

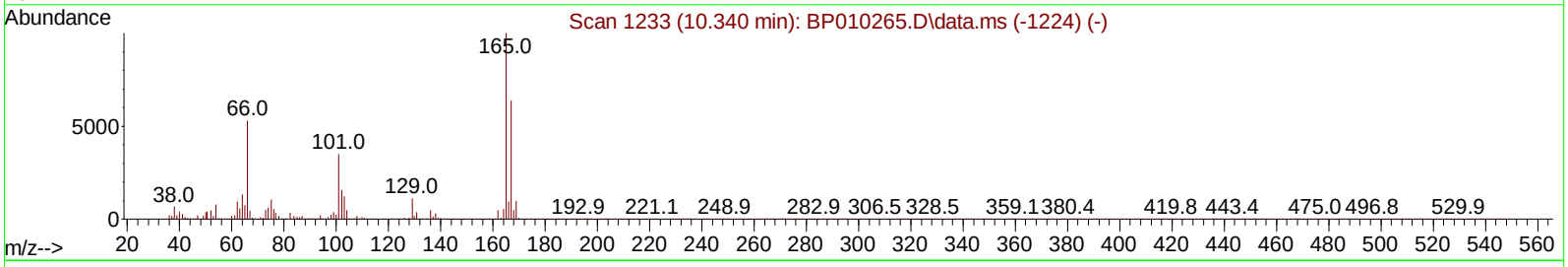

TIC: BP010270.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.463min (+ 0.123) 0.03 ng/ul

response 213

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	53.52
101.00	49.80	42.95
0.00	0.00	0.00

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Misc :
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ClientSampleId :
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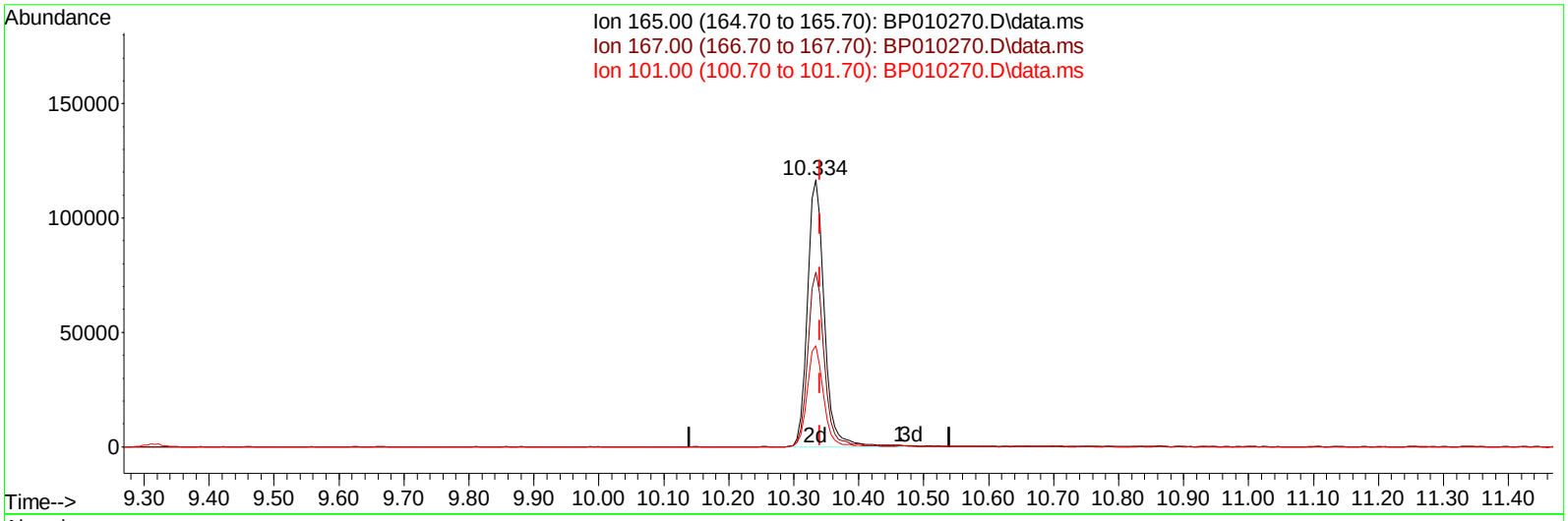

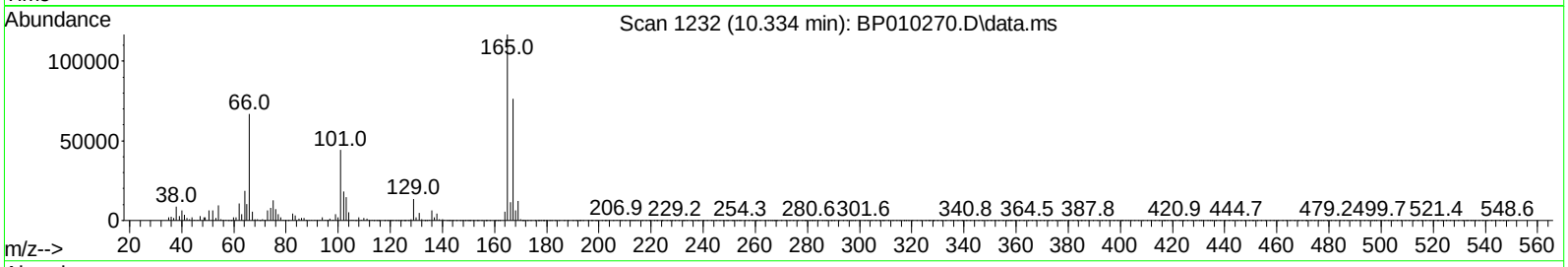

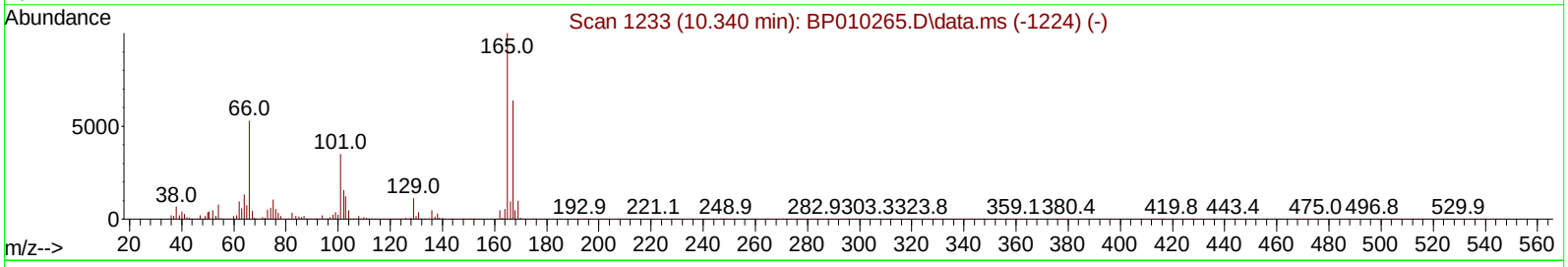

TIC: BP010270.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.334min (-0.006) 28.88 ng/ul m

response 212791

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	65.54
101.00	49.80	37.93#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051022\
Data File : BP010270.D
Acq On : 10 May 2022 17:59
Operator : CG/JU
Sample : N2678-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
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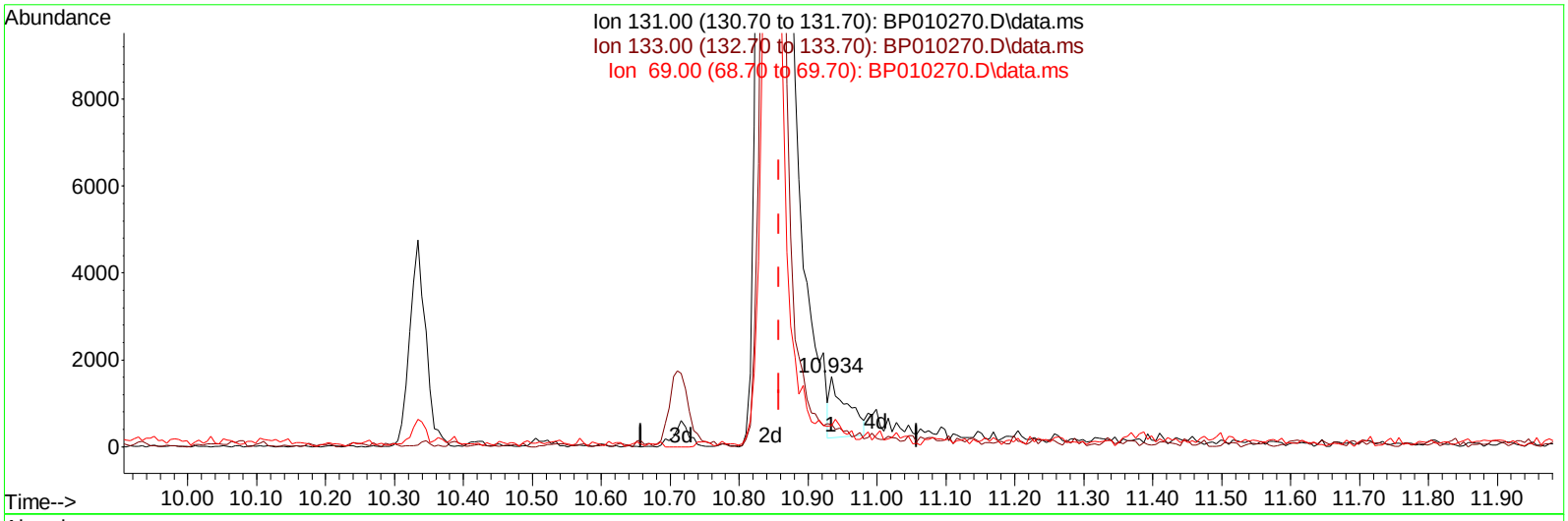
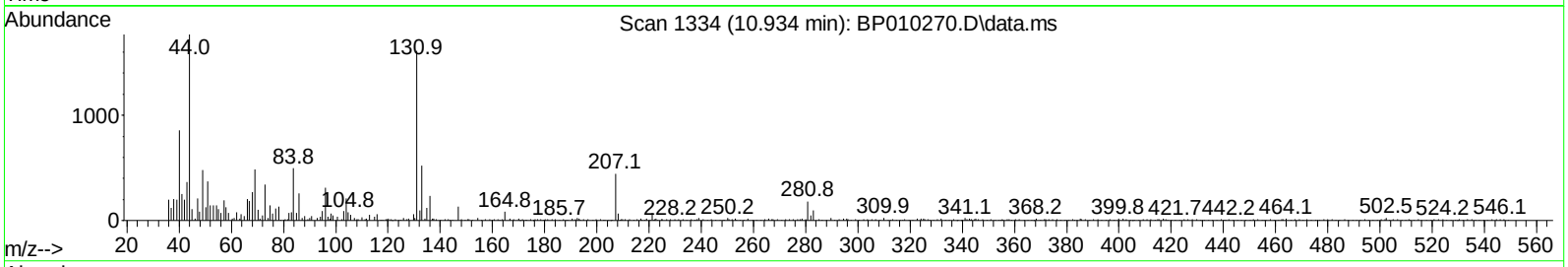
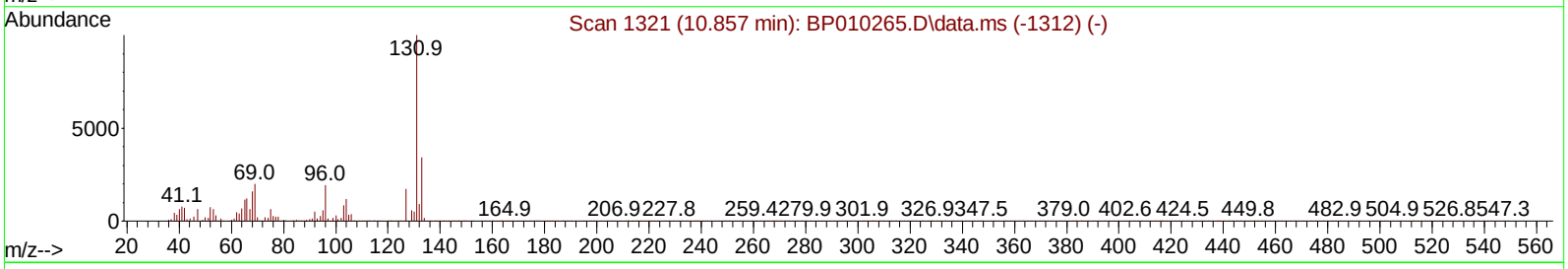
Instrument :
BNA_P
ClientSampleId :
BH0C1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/11/2022
Supervised By :mohammad ahmed 05/12/2022







TIC: BP010270.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.934min (+ 0.076) 0.22 ng/ul

response 2395

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	32.42
69.00	26.10	30.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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Instrument :
BNA_P
ClientSampleId :
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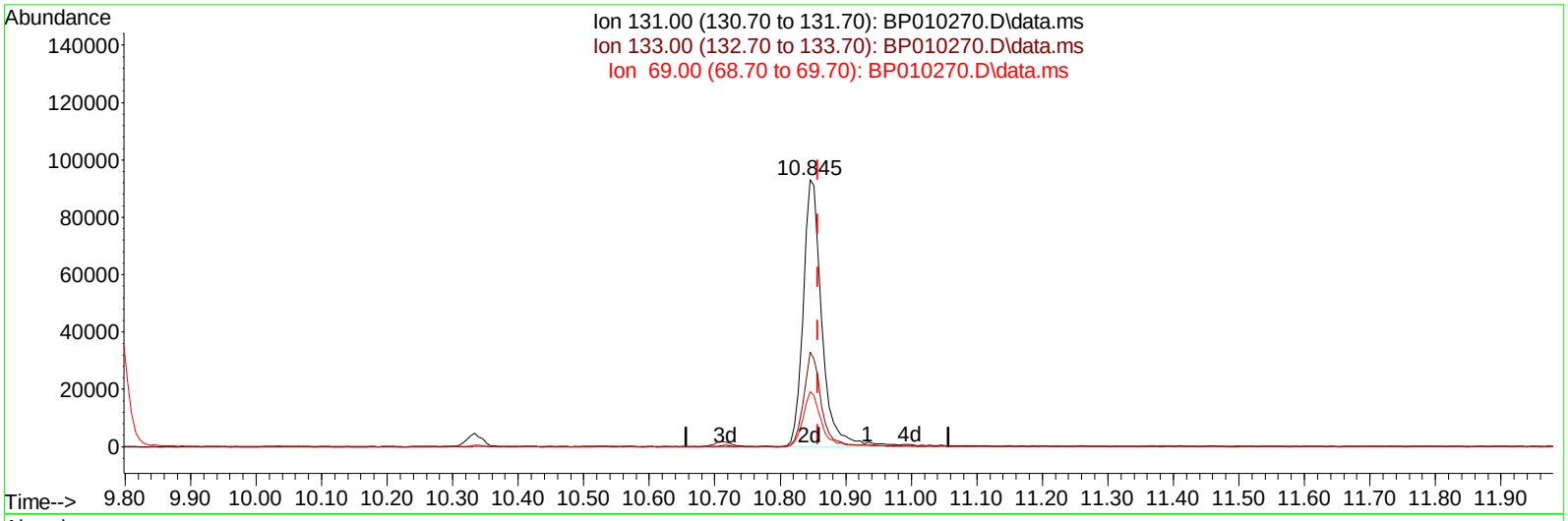
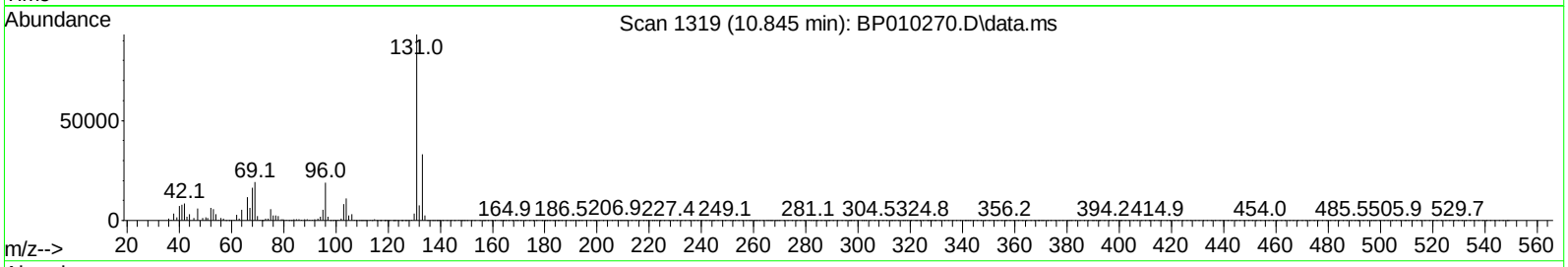
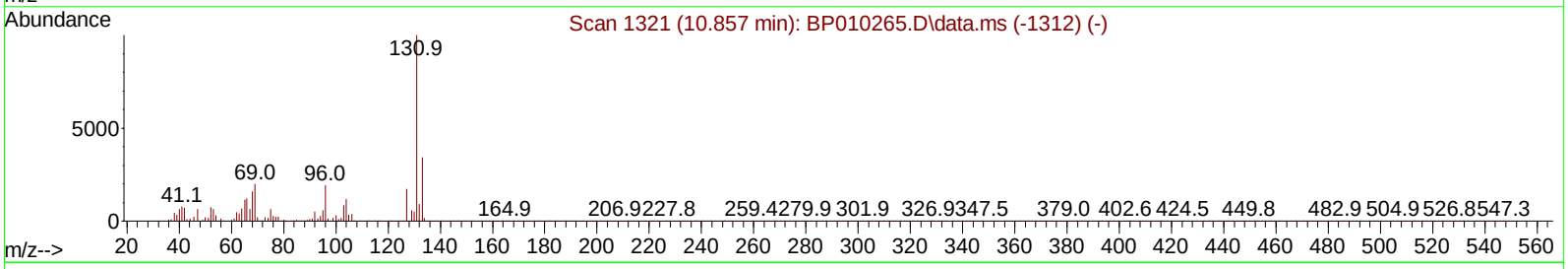
Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/11/2022
Supervised By :mohammad ahmed 05/12/2022







TIC: BP010270.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.845min (-0.012) 17.08 ng/ul m

response 185828

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	35.47#
69.00	26.10	20.60#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051022\
Data File : BP010270.D
Acq On : 10 May 2022 17:59
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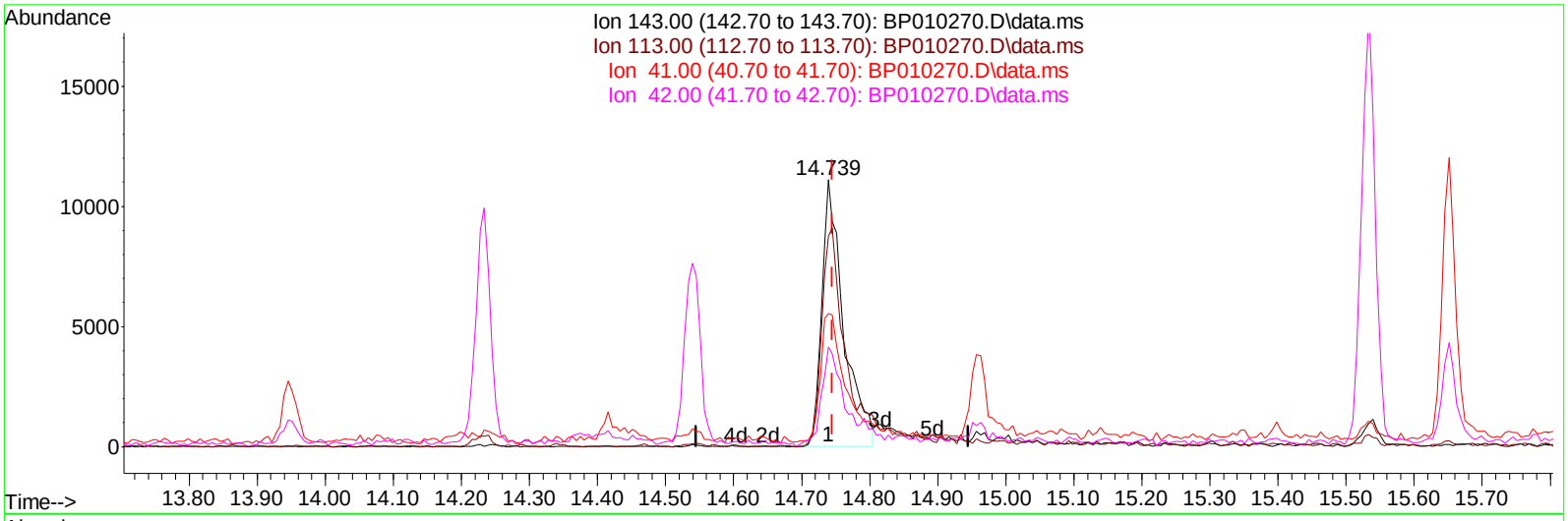
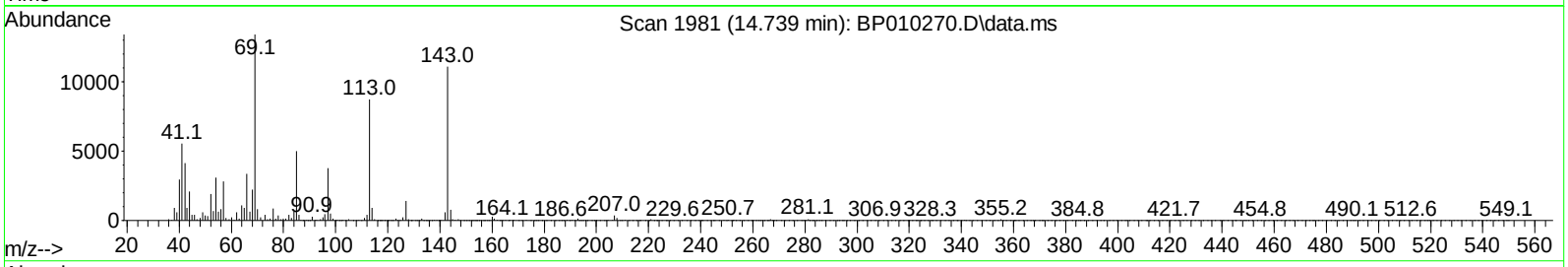
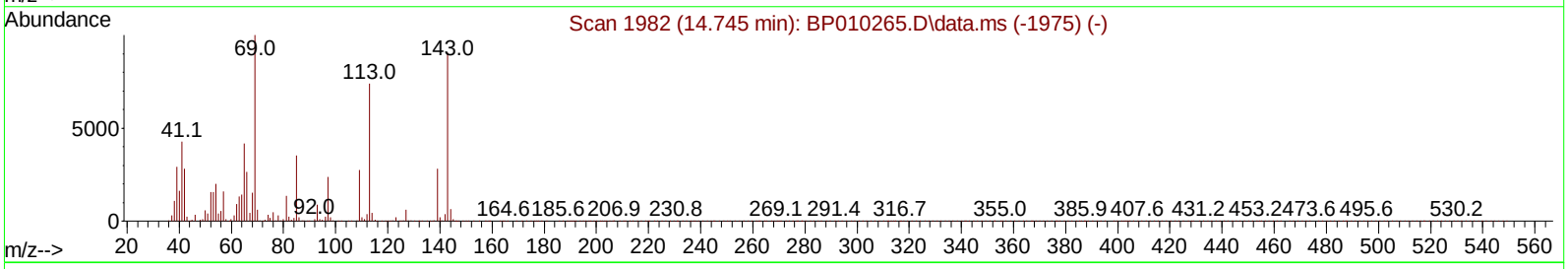
Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/11/2022
Supervised By :mohammad ahmed 05/12/2022







TIC: BP010270.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.006) 4.97 ng/ul

response 24649

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.78#
41.00	23.20	49.92#
42.00	18.00	37.42#

Quantitation Report (Qedit)

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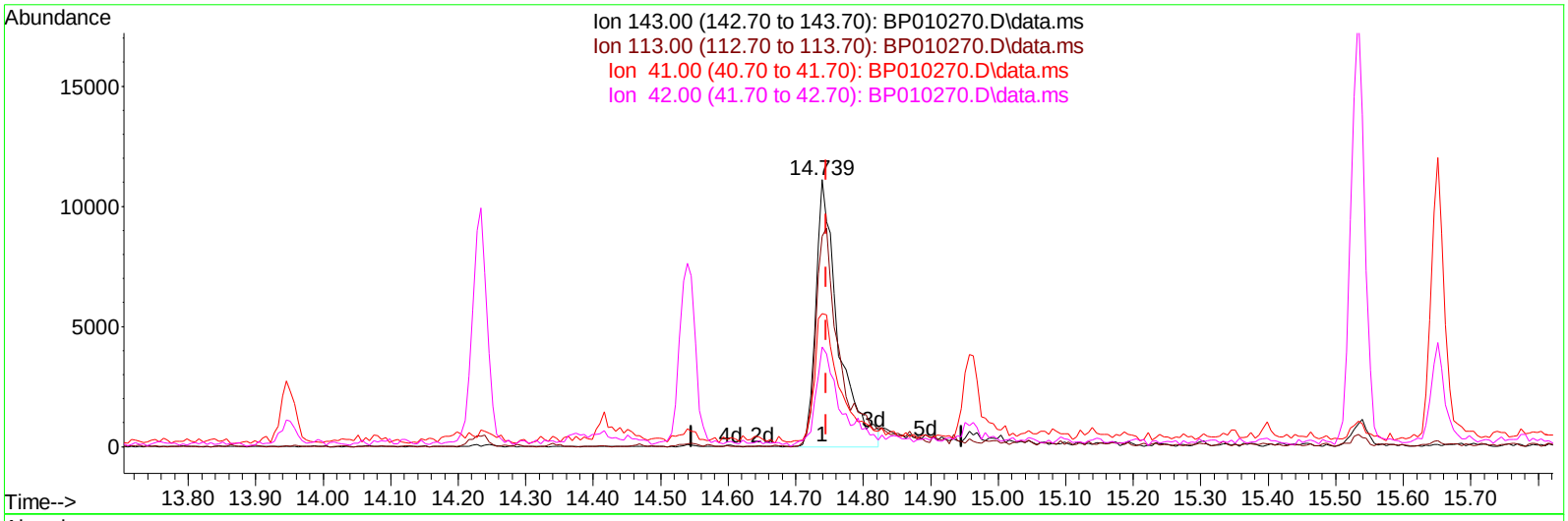

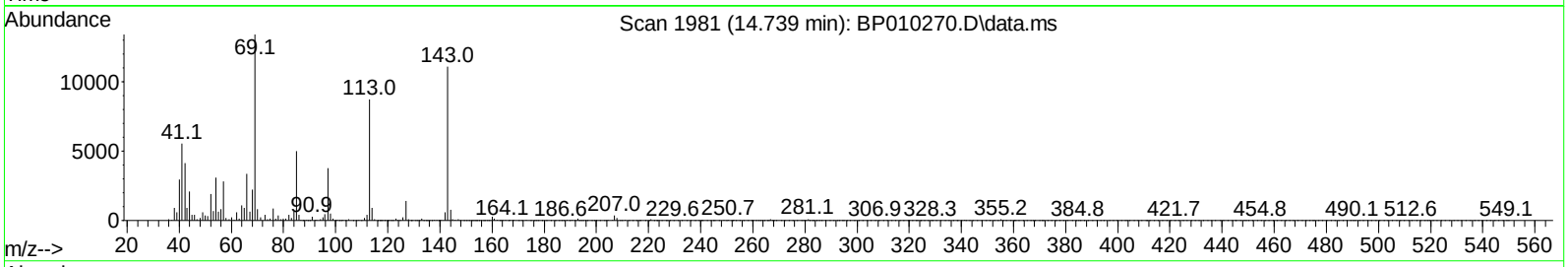

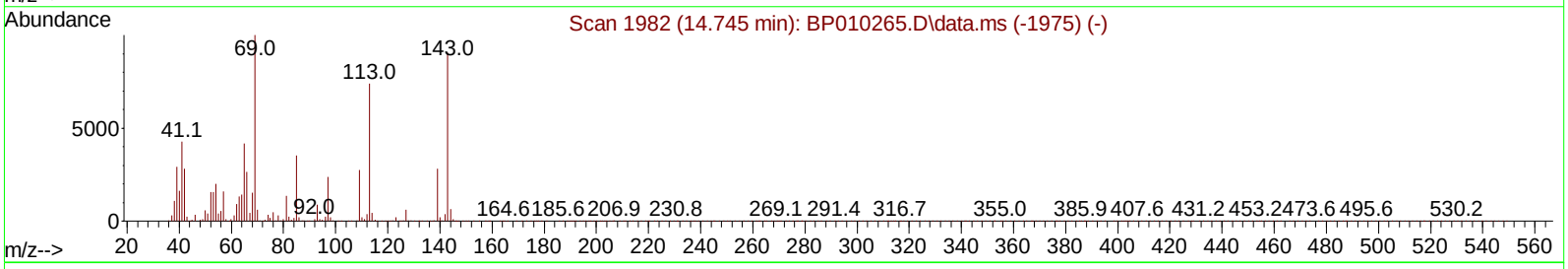

TIC: BP010270.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.006) 5.18 ng/ul m

response 25659

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.78#
41.00	23.20	49.92#
42.00	18.00	37.42#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	102858	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.710	136	442866	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.539	164	324282	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.292	188	728562	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.374	240	741049	20.000	ng/ul	#-0.01
88) Perylene-d12	23.810	264	781490	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	13885m	5.735	ng/uL	0.00
4) Pyridine-d5	3.781	84	46285	6.870	ng/ul	0.01
7) Phenol-d5	7.075	99	62966	6.591	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	201930	34.168	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	181747	25.832	ng/ul	0.00
15) 4-Methylphenol-d8	8.616	113	125425	15.969	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	123693	34.834	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	127531	31.499	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	212791m	28.876	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	185828m	17.081	ng/ul	-0.01
46) Dimethylphthalate-d6	13.951	166	957475	37.062	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	992612	32.226	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	25659m	5.178	ng/ul	0.00
60) Fluorene-d10	15.533	176	818019	37.511	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.651	200	155109	31.089	ng/ul	-0.01
73) Anthracene-d10	17.392	188	1403212	39.330	ng/ul	-0.01
81) Pyrene-d10	19.621	212	1756559	42.894	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.657	264	1754610	42.100	ng/ul	-0.01
Target Compounds						Qvalue

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

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