Quantitation Report (Qedit)

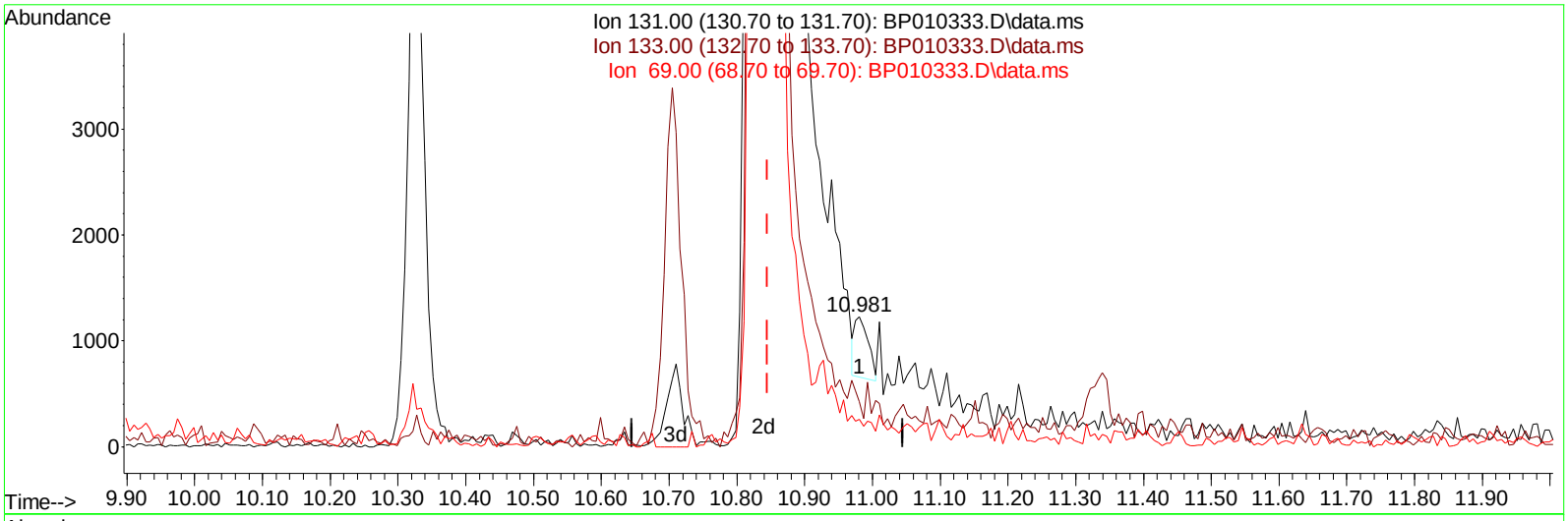
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010333.D
Acq On : 12 May 2022 20:50
Operator : CG/JU
Sample : N2746-06
Misc :
ALS Vial : 57 Sample Multiplier: 1

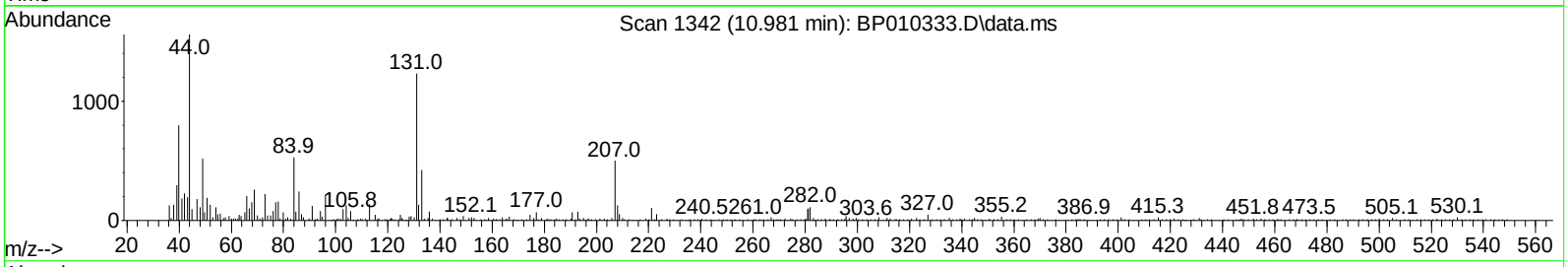
Quant Time: May 13 00:03:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 12 01:10:15 2022
Response via : Initial Calibration

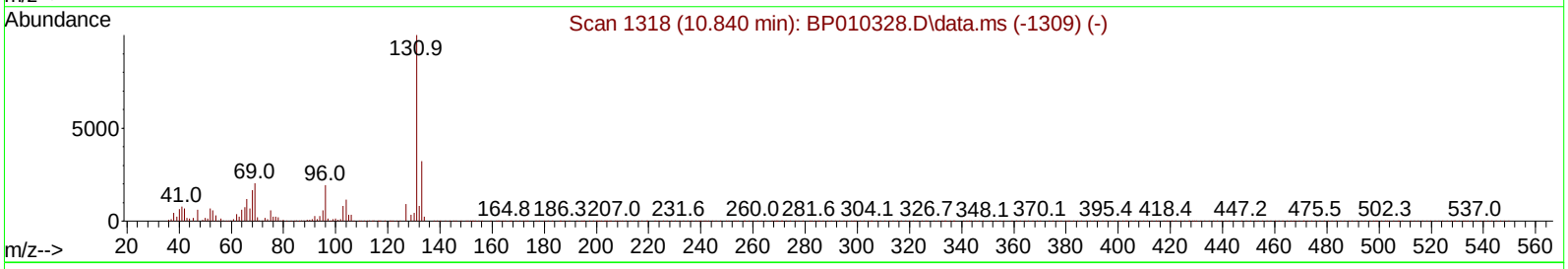
Instrument :
BNA_P
ClientSampleId :
DBSL2

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/16/2022
Supervised By :Yogesh Patel 05/17/2022







TIC: BP010333.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.981min (+ 0.135) 0.05 ng/ul

response 793

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	34.42
69.00	26.10	21.07
0.00	0.00	0.00

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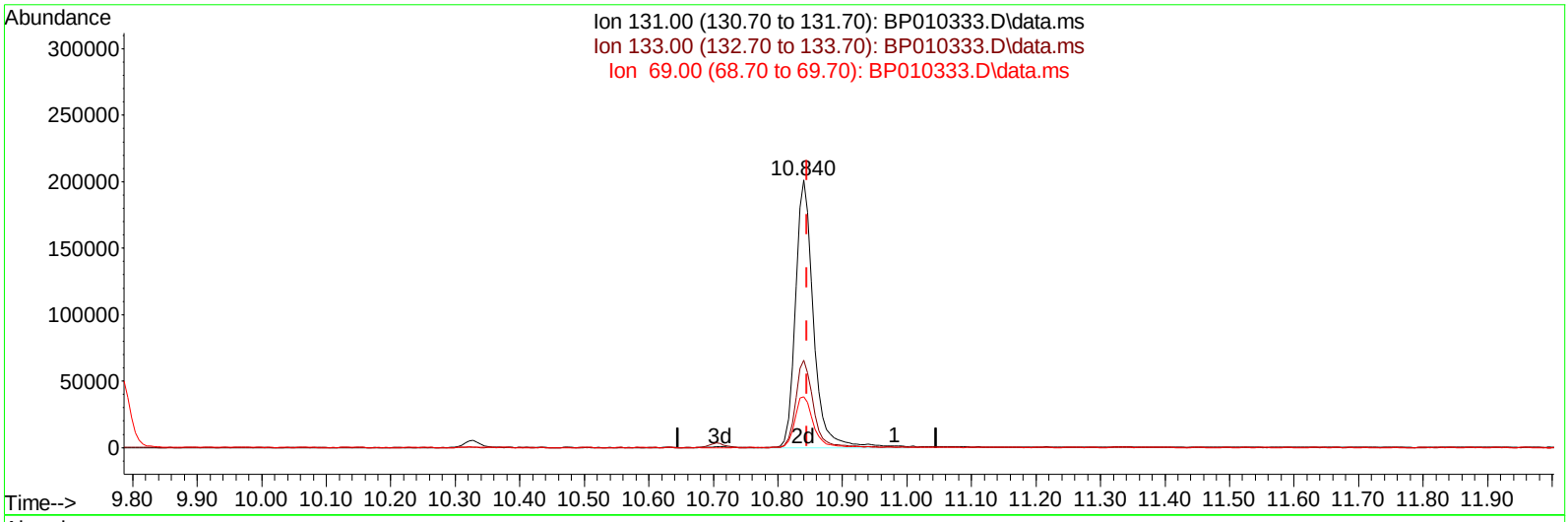

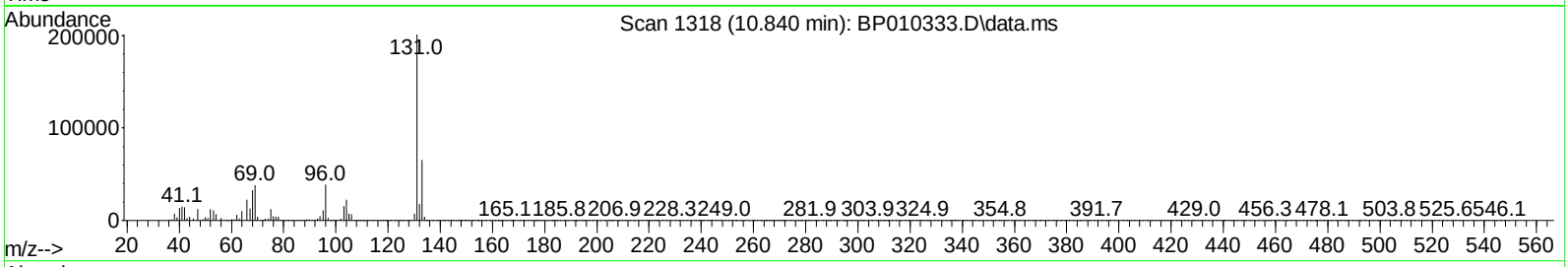

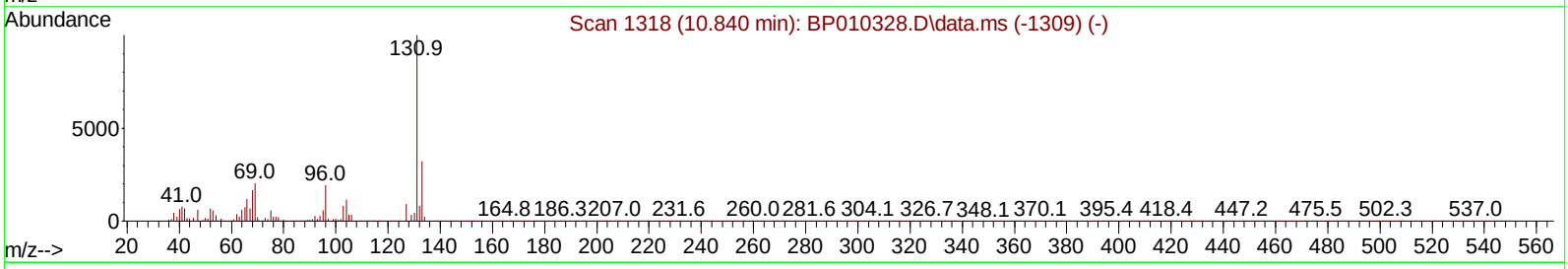

TIC: BP010333.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.840min (-0.006) 22.99 ng/ul m

response 391589

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	32.64
69.00	26.10	18.93#
0.00	0.00	0.00

Quantitation Report (Qedit)

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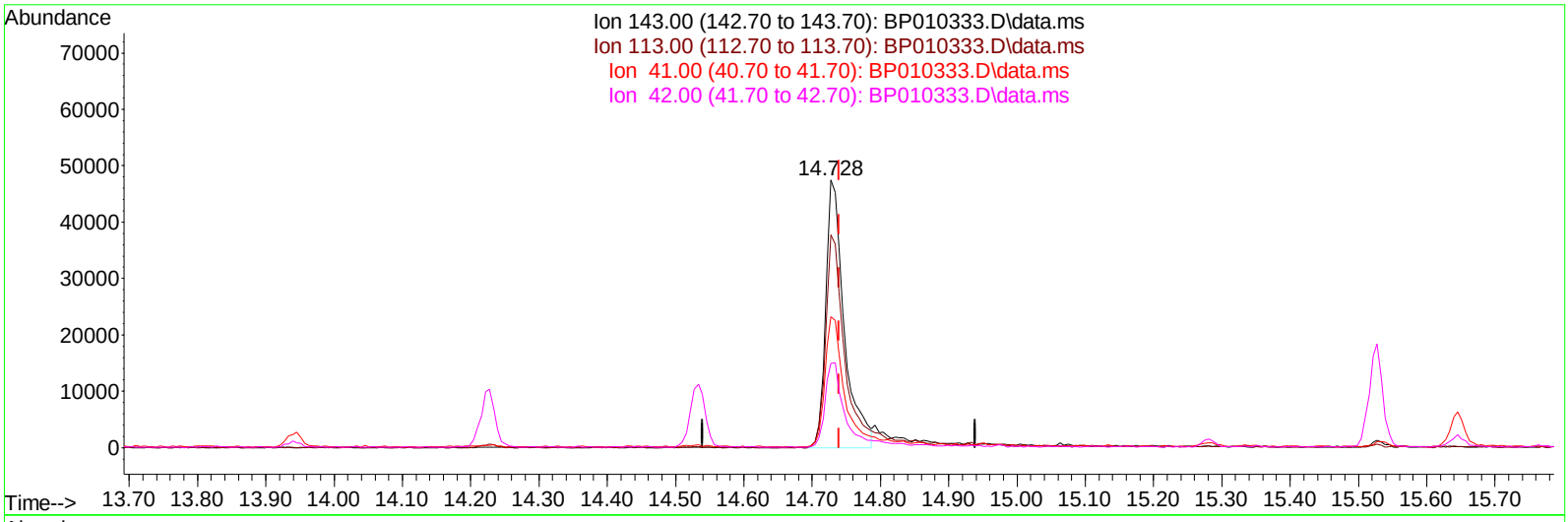

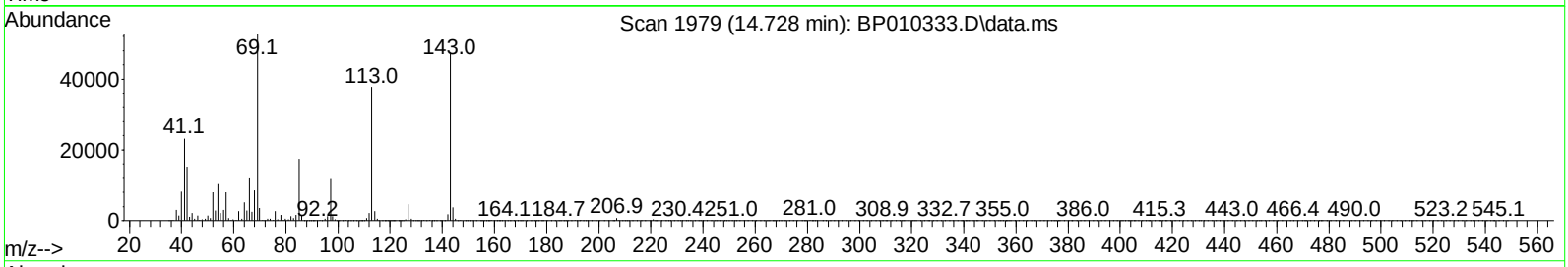

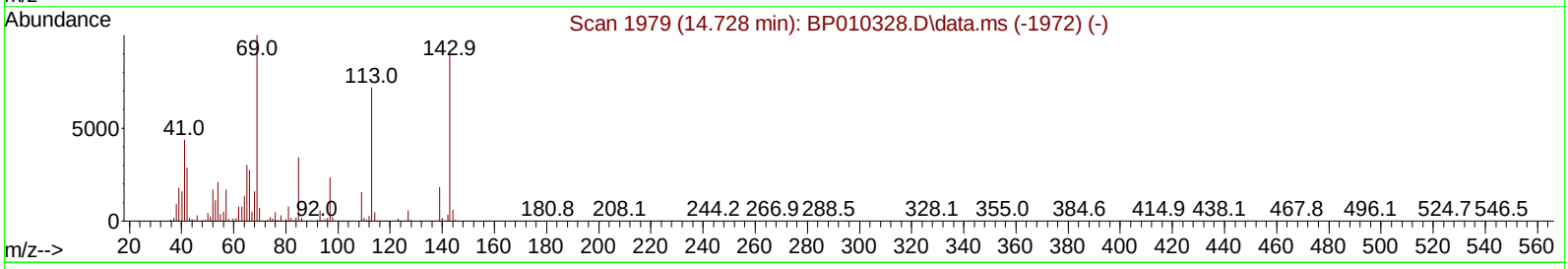

TIC: BP010333.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.012) 11.91 ng/ul

response 89739

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.66#
41.00	23.20	49.05#
42.00	18.00	31.56#

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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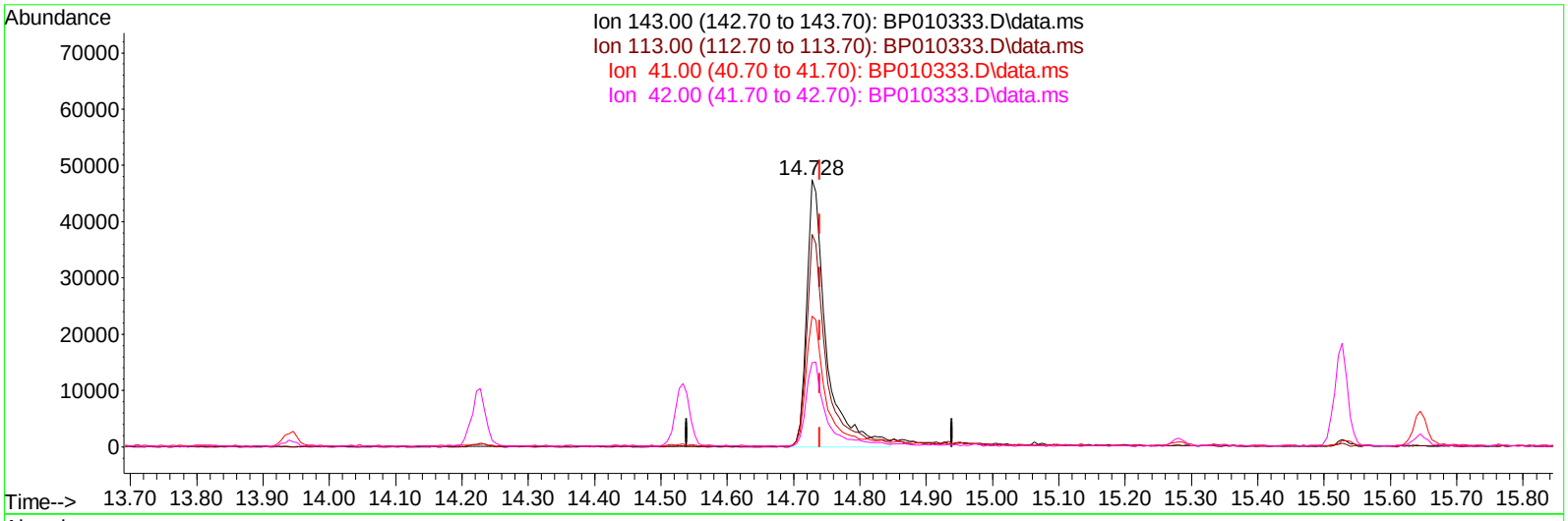

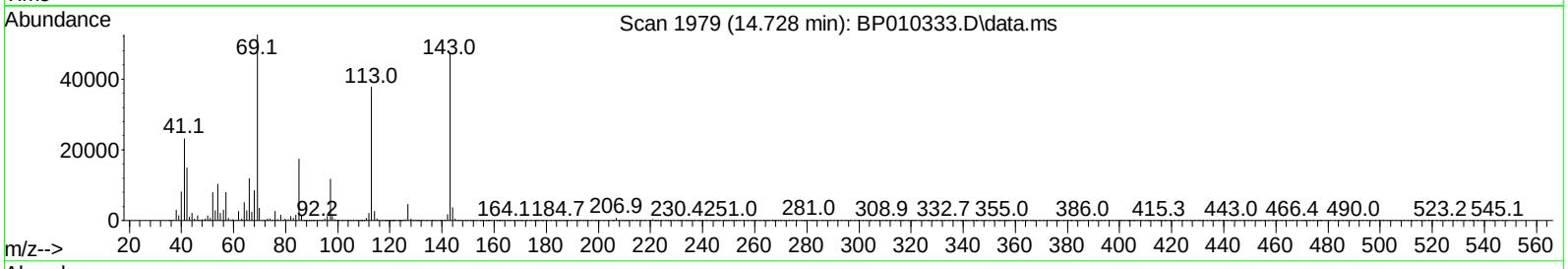

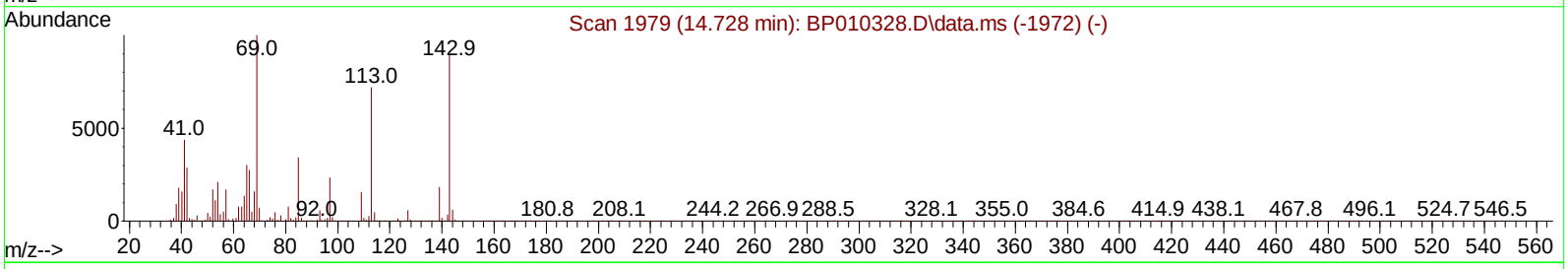

TIC: BP010333.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.012) 12.89 ng/ul m

response 97094

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.66#
41.00	23.20	49.05#
42.00	18.00	31.56#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 05/17/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	156720	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	693310	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.533	164	492994	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1080684	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	1052849	20.000	ng/ul	#-0.01
88) Perylene-d12	23.792	264	918382	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	15067	4.084	ng/uL	0.00
4) Pyridine-d5	3.763	84	222352	21.659	ng/ul	0.00
7) Phenol-d5	7.069	99	312271	21.453	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	219373	24.362	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	250562	23.373	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	275392	23.013	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	137502	24.735	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	150885	23.805	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	266193	23.074	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	391589m	22.992	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	992608	25.273	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1106221	23.624	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	97094m	12.888	ng/ul	-0.01
60) Fluorene-d10	15.528	176	848441	25.592	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	92026	12.435	ng/ul	-0.01
73) Anthracene-d10	17.386	188	1316436	24.875	ng/ul	0.00
81) Pyrene-d10	19.616	212	1623341	27.901	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264	1316792	26.886	ng/ul	-0.02
Target Compounds						Qvalue

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

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