Quantitation Report (Qedit)

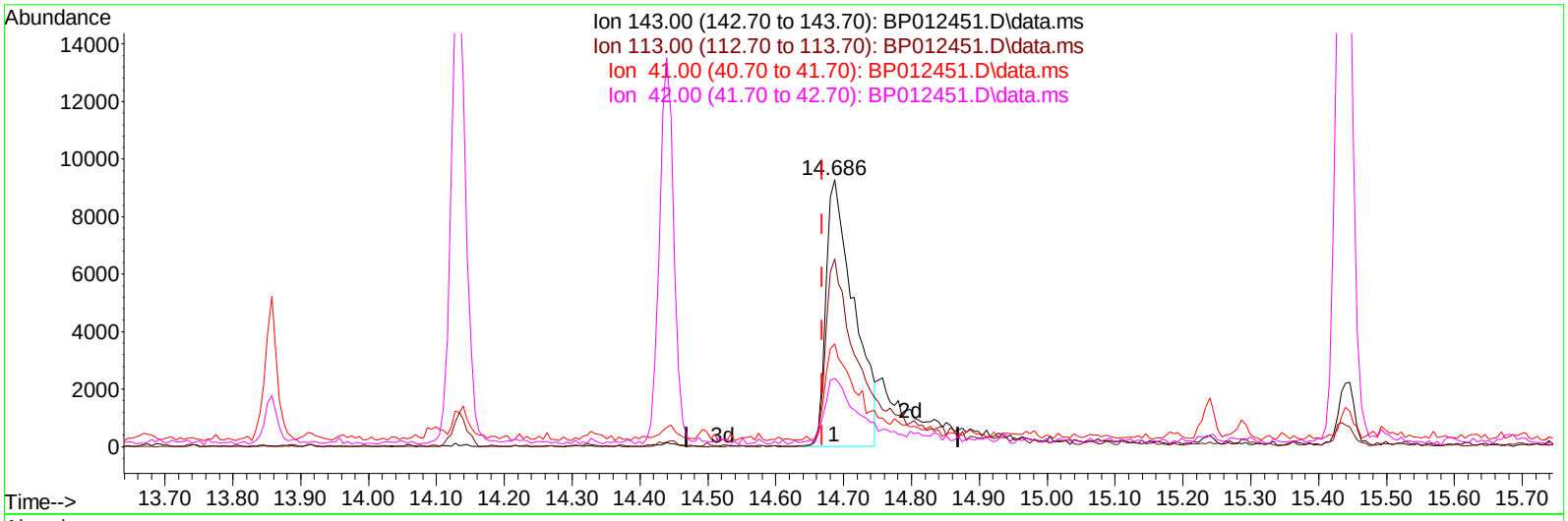
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012451.D
Acq On : 02 Nov 2022 13:37
Operator : CG/JU
Sample : N5180-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

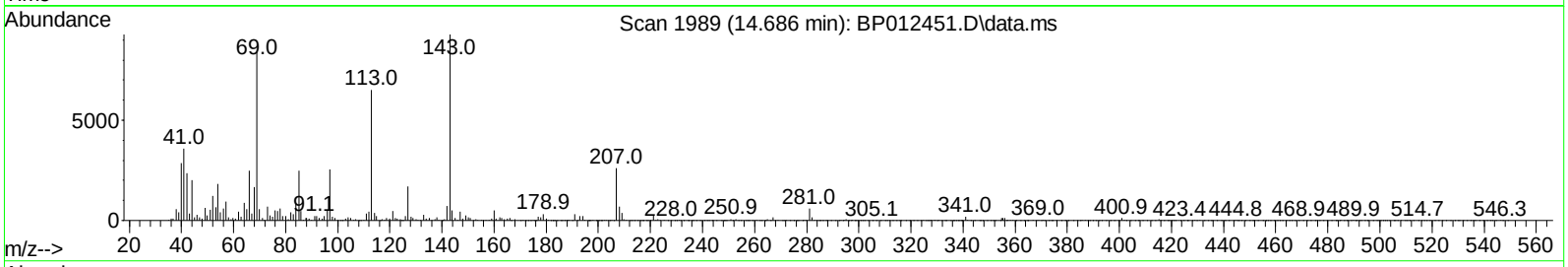
Quant Time: Nov 02 23:14:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 01 01:03:58 2022
Response via : Initial Calibration

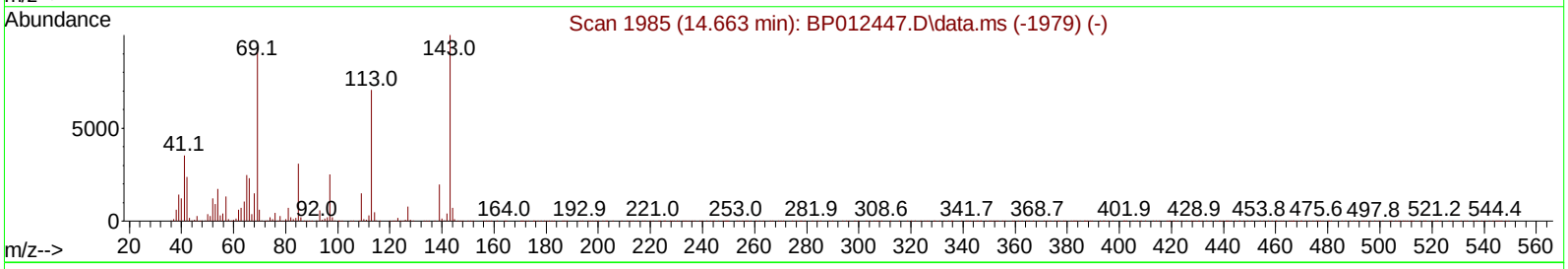
Instrument :
BNA_P
ClientSampleId :
C0KX3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/03/2022
Supervised By :mohammad ahmed 11/03/2022







TIC: BP012451.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.017) 2.51 ng/ul

response 26505

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	70.30
41.00	34.50	38.60
42.00	23.60	25.58

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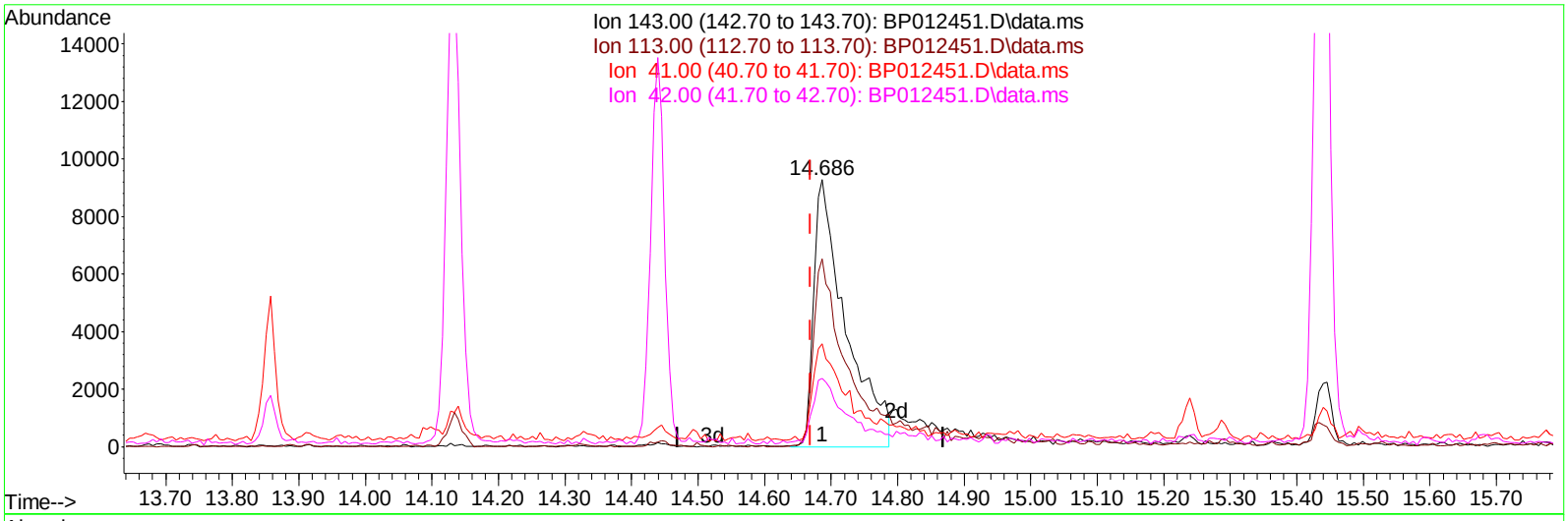

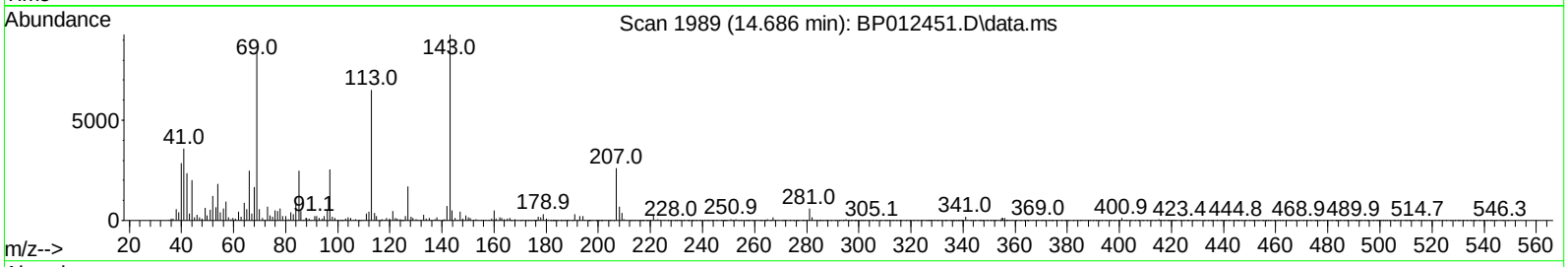

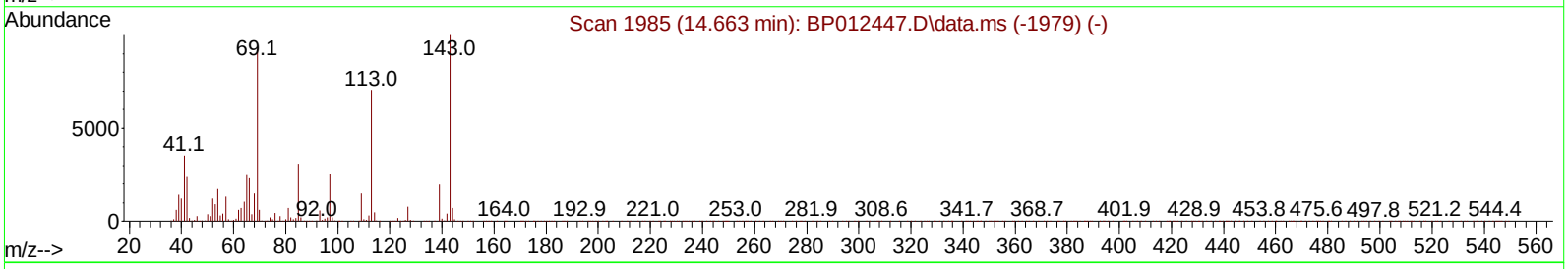

TIC: BP012451.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (+ 0.017) 2.92 ng/ul m

response 30913

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	70.30
41.00	34.50	38.60
42.00	23.60	25.58

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ClientSampleId :
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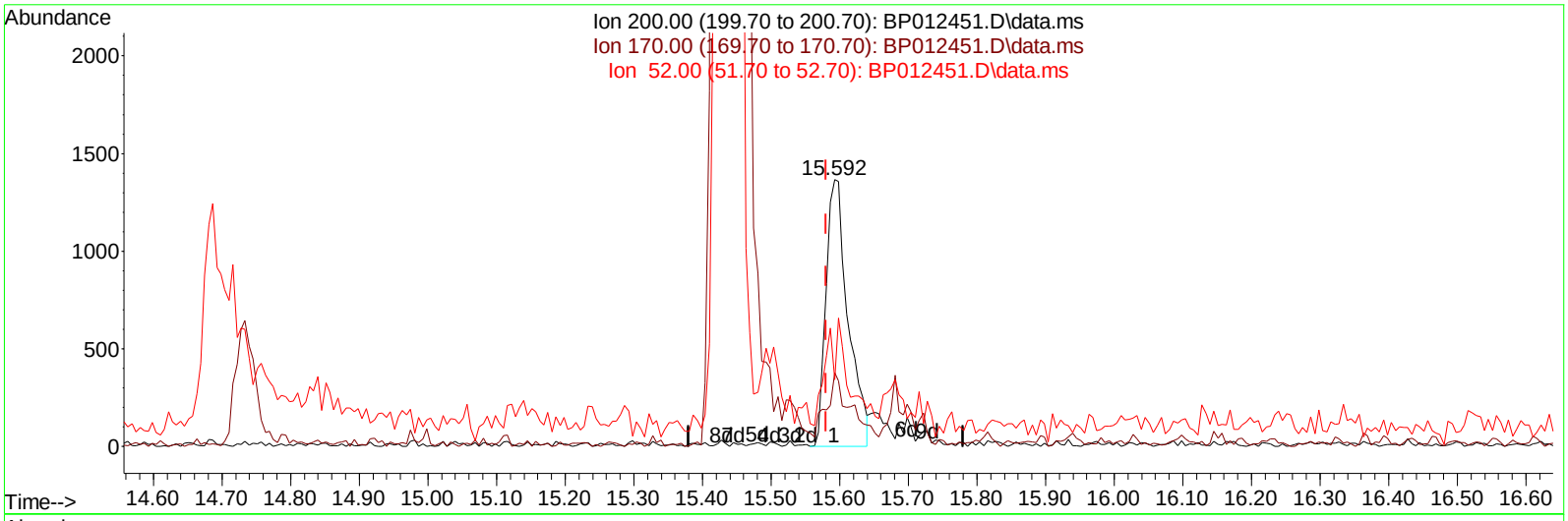

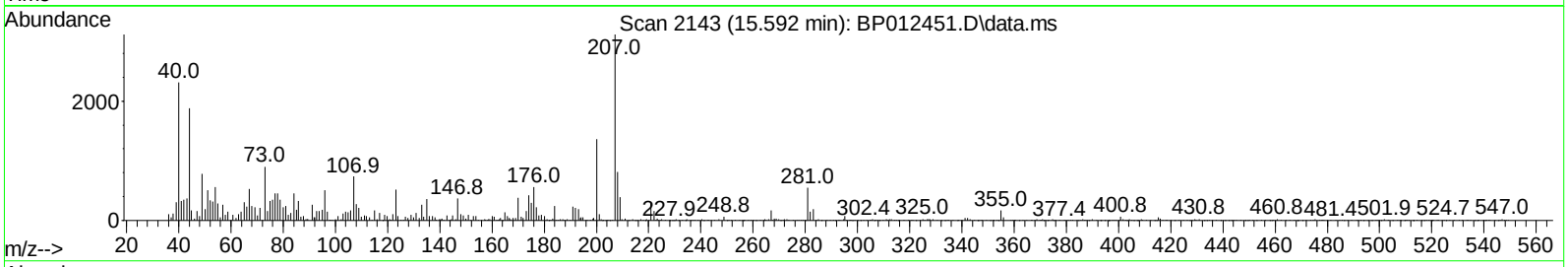

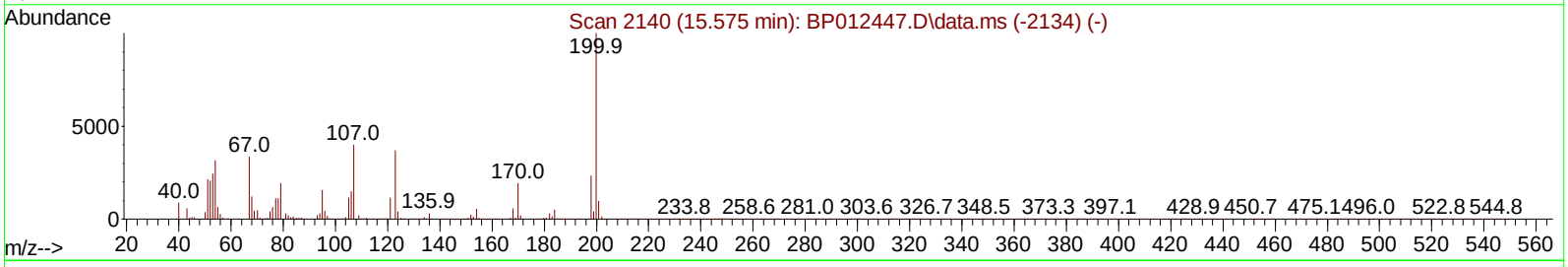

TIC: BP012451.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.592min (+ 0.012) 0.32 ng/ul

response 3059

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.90	27.75#
52.00	37.70	25.11#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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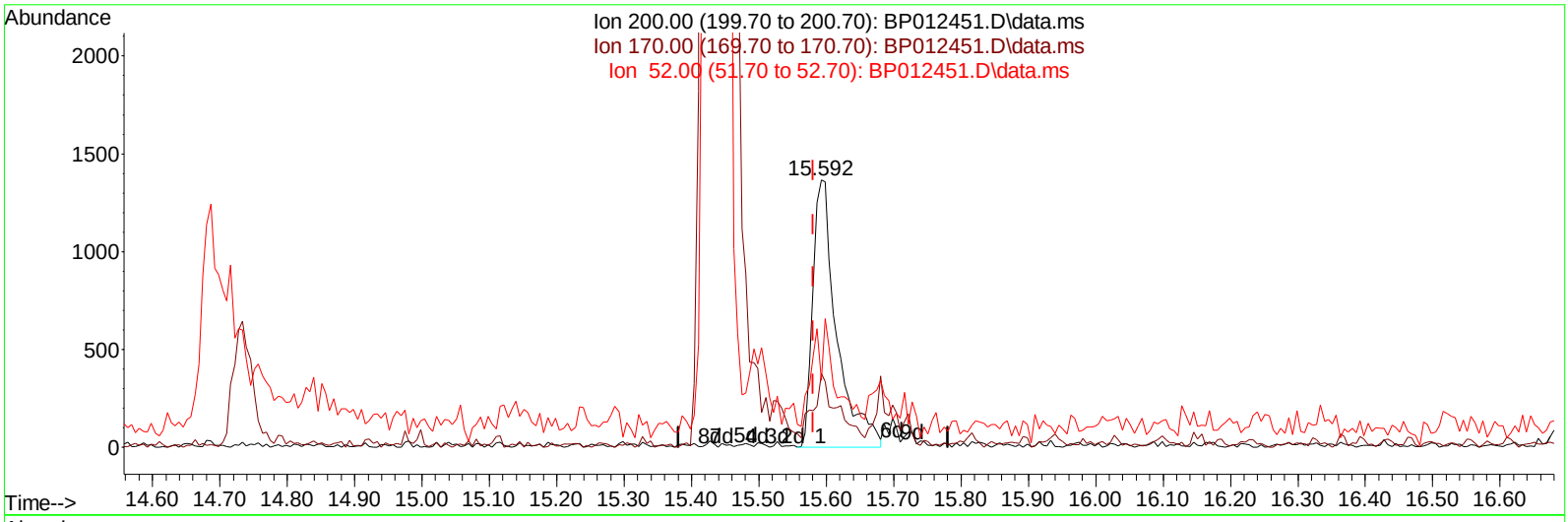

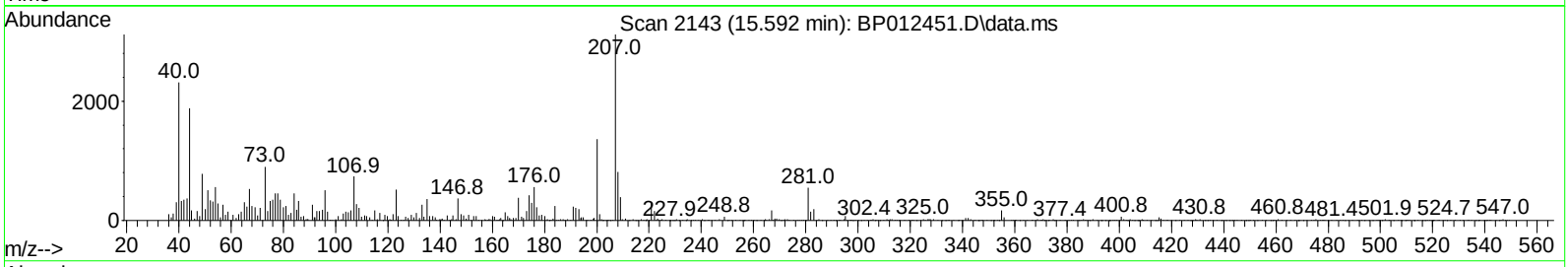

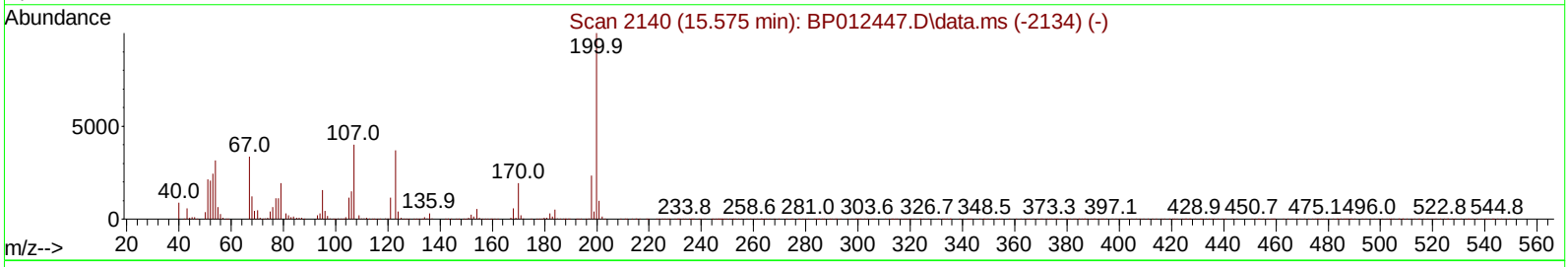

TIC: BP012451.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.592min (+ 0.012) 0.35 ng/ul m

response 3366

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.90	27.75#
52.00	37.70	25.11#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :mohammad ahmed 11/03/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	262785	20.000	ng/ul	0.00
20) Naphthalene-d8	10.587	136	1189023	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.439	164	816579	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1828721	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	1698662	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1523420	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	27170	4.163	ng/uL	0.00
4) Pyridine-d5	3.664	84	150767	8.042	ng/ul	0.00
7) Phenol-d5	6.963	99	229715	9.746	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	491055	34.900	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	339067	19.699	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	327564	17.572	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	330030	42.802	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	58547	7.316	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	346267	19.965	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	766540	29.683	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2228980	39.875	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	2475731	38.756	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	30913m	2.922	ng/ul	0.02
60) Fluorene-d10	15.439	176	2013929	42.080	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	3366m	0.354	ng/ul	0.01
73) Anthracene-d10	17.304	188	3540262	44.930	ng/ul	0.00
81) Pyrene-d10	19.545	212	4013194	47.057	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.533	264	3364508	45.200	ng/ul	-0.01
Target Compounds						Qvalue

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

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