Quantitation Report (Qedit)

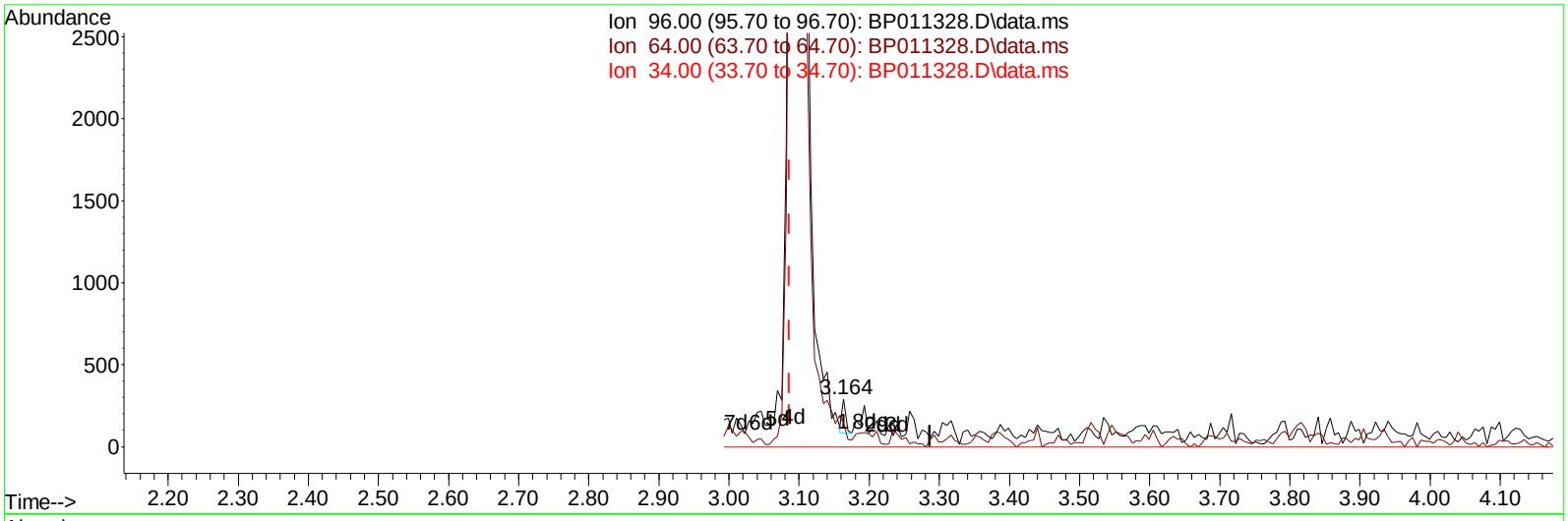
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011328.D
Acq On : 09 Aug 2022 14:20
Operator : CG/JU
Sample : N4026-21
Misc :
ALS Vial : 10 Sample Multiplier: 1

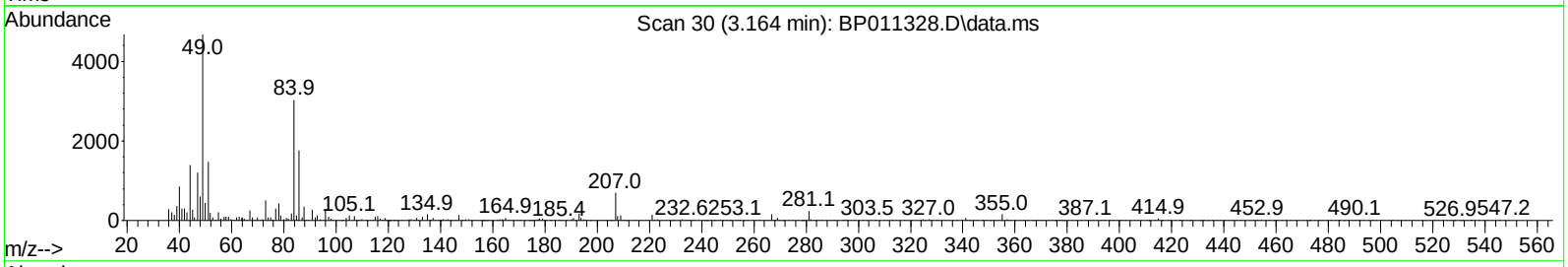
Quant Time: Aug 09 23:11:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

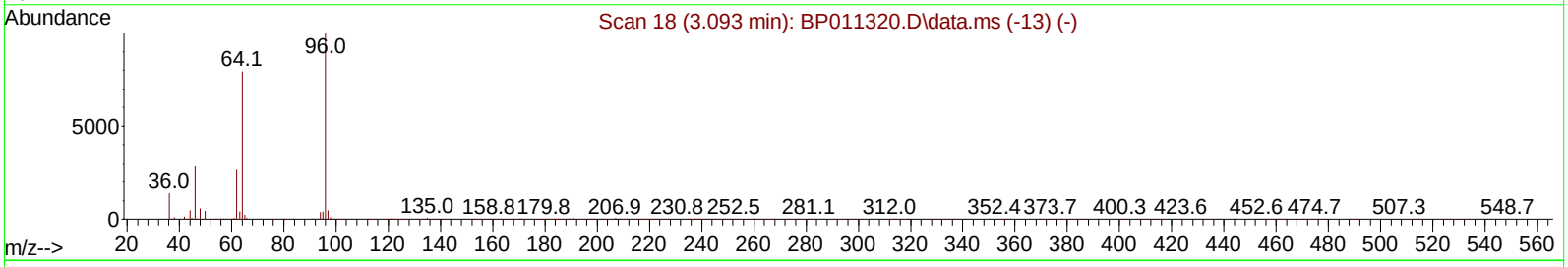
Instrument :
BNA_P
ClientSampleId :
GBJJ6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/10/2022







TIC: BP011328.D\data.ms

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

3.164min (+ 0.077) 0.02 ng/uL

response 79

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	50.17
34.00	0.00	0.00
0.00	0.00	0.00

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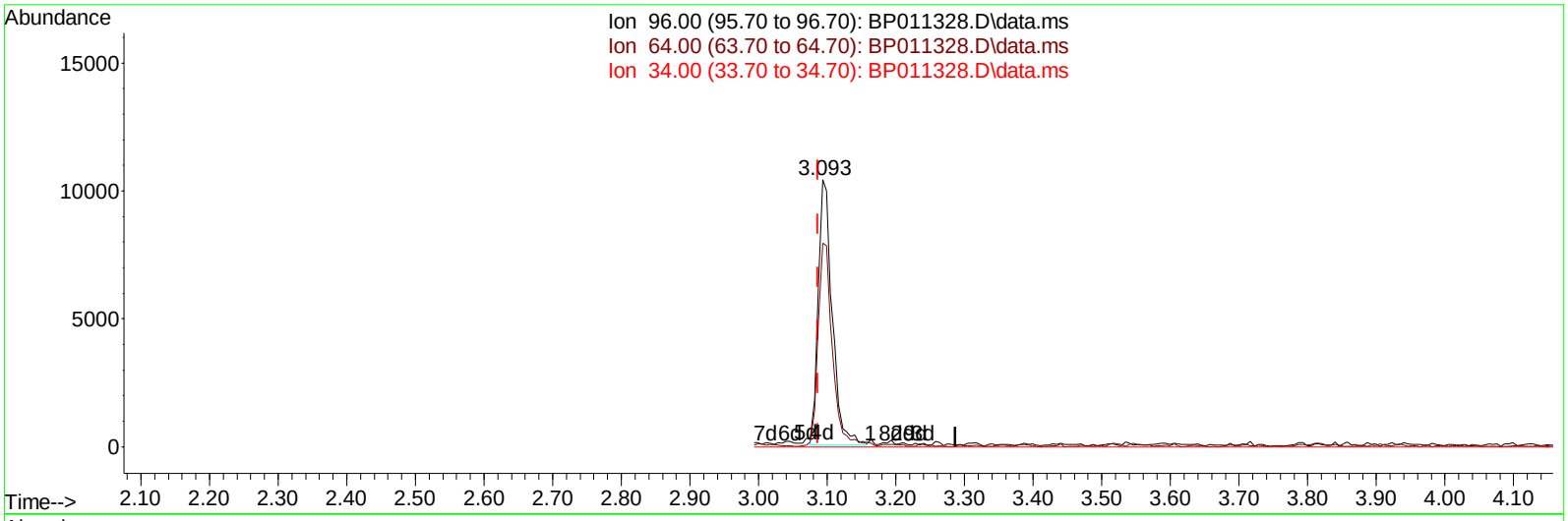

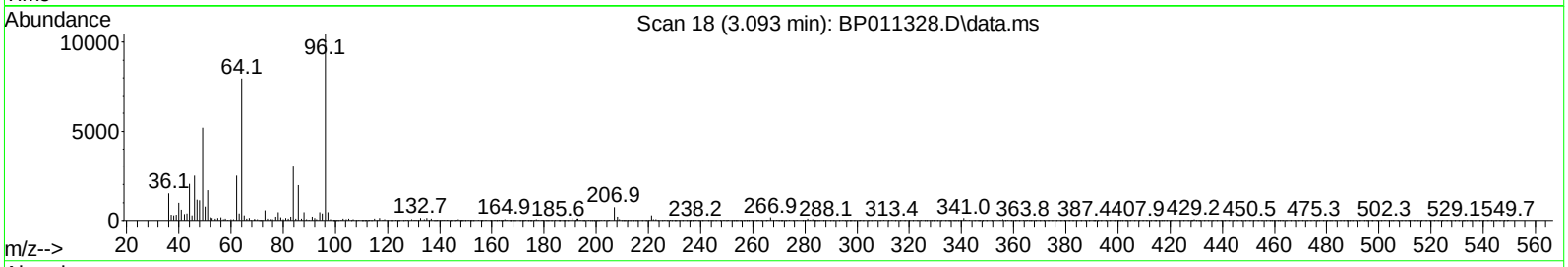

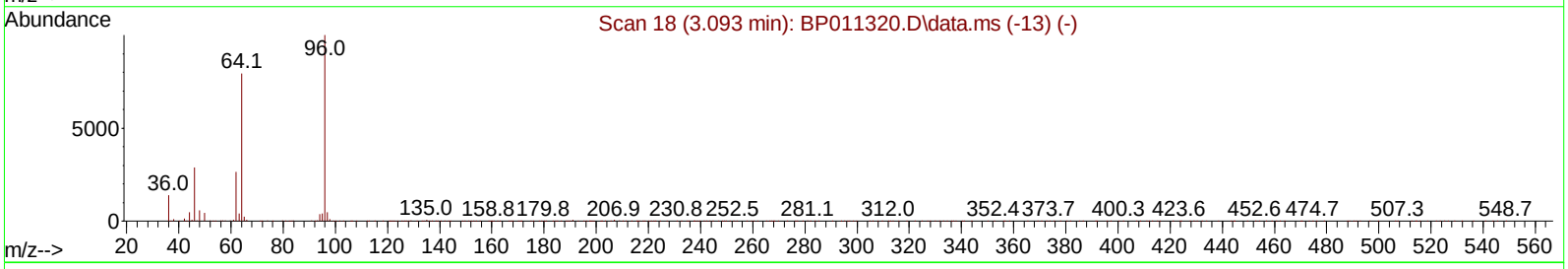

TIC: BP011328.D\data.ms

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

3.093min (+ 0.006) 3.73 ng/uL m

response 14785

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	76.21#
34.00	0.00	0.00
0.00	0.00	0.00

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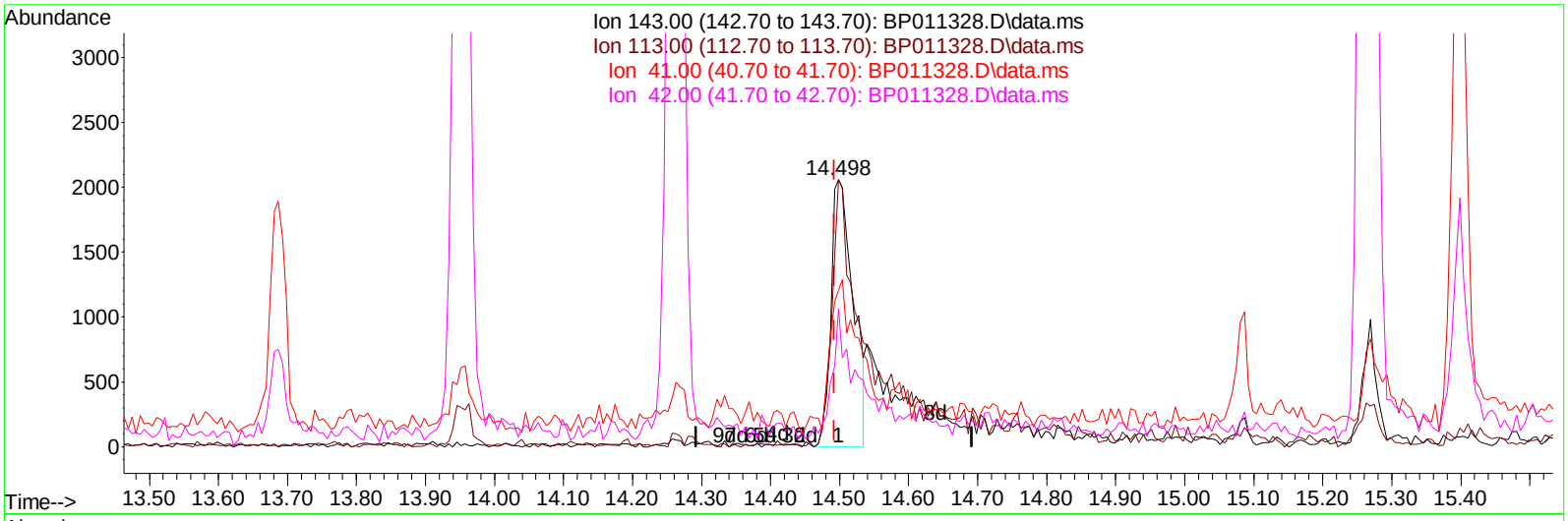

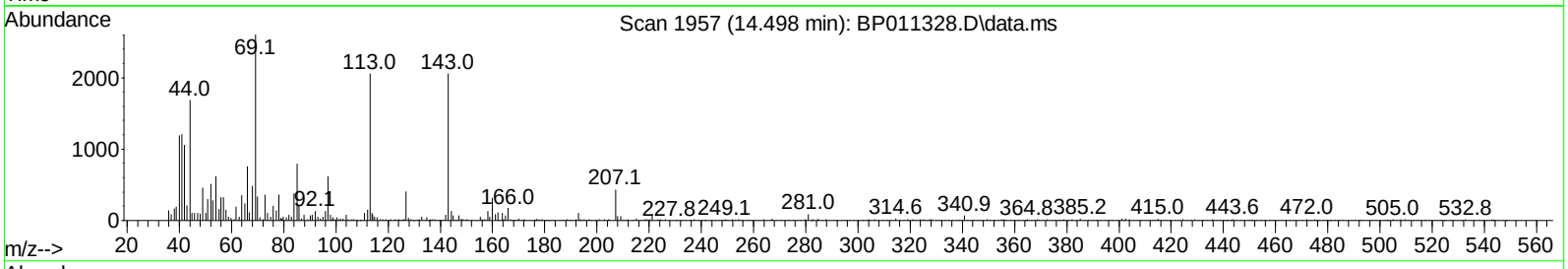

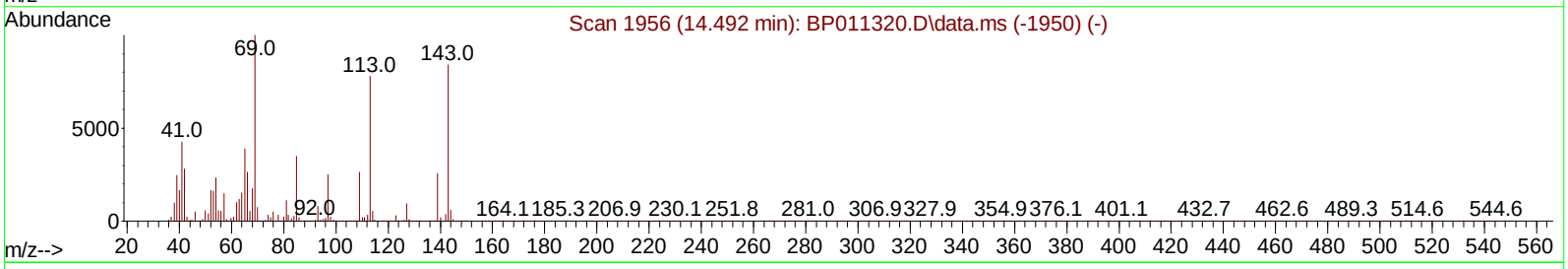

TIC: BP011328.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 1.25 ng/ul

response 4594

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	100.00#
41.00	54.40	58.65
42.00	35.50	51.70#

Quantitation Report (Qedit)

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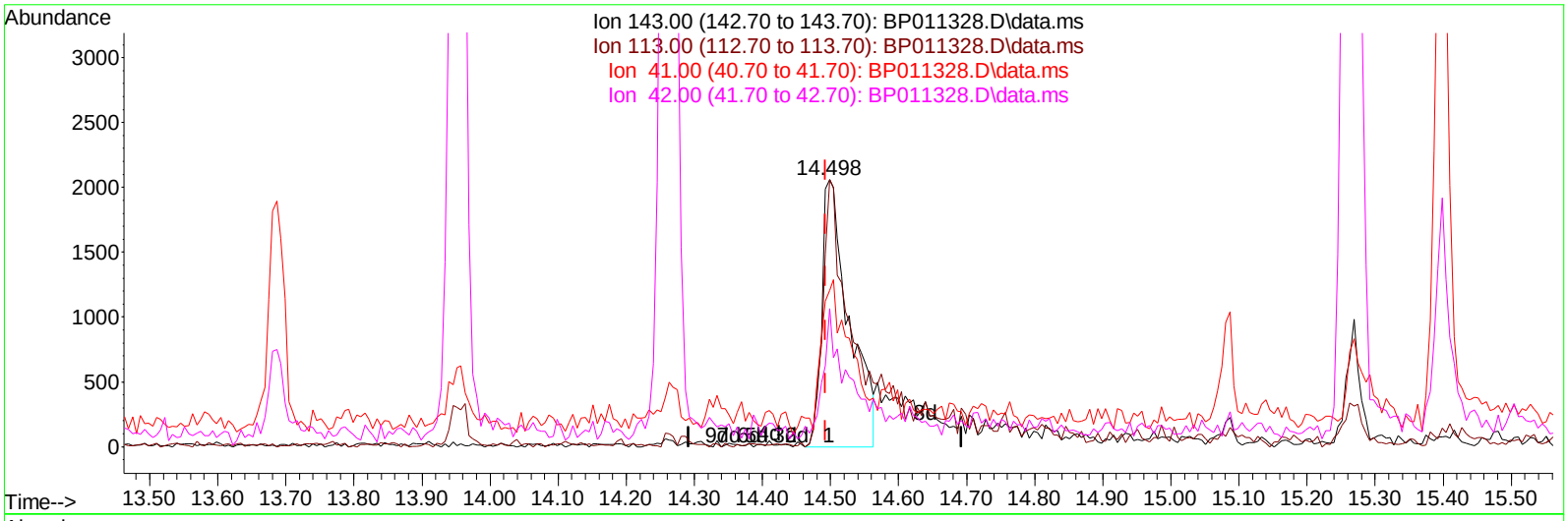

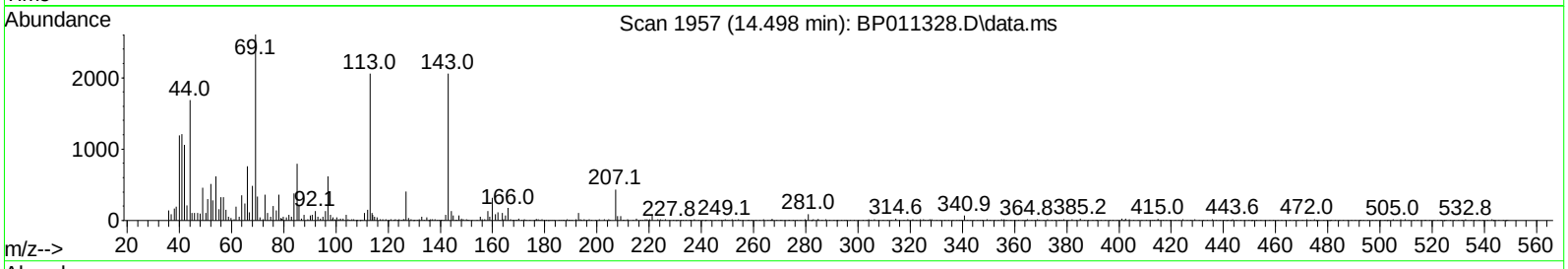

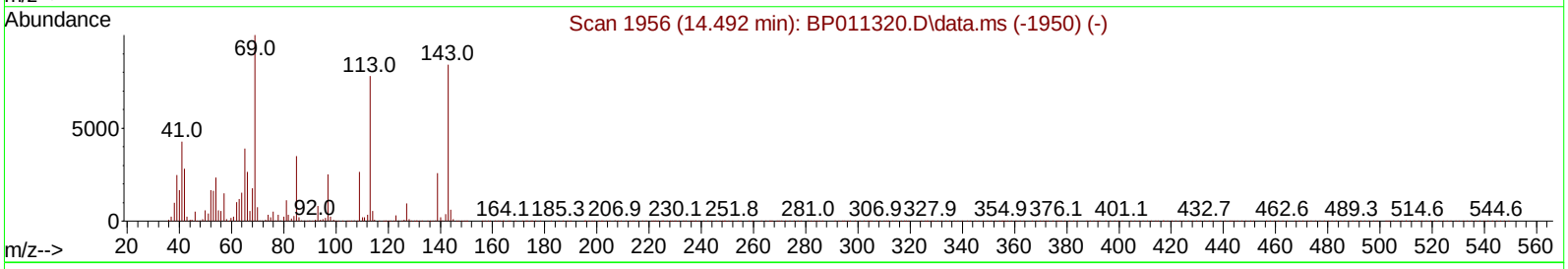

TIC: BP011328.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 1.55 ng/ul m

response 5691

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	100.00#
41.00	54.40	58.65
42.00	35.50	51.70#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	121603	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	512765	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	289392	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	578632	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	532351	20.000	ng/ul	0.00
88) Perylene-d12	23.474	264	527786	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	14785m	3.727	ng/uL	0.00
4) Pyridine-d5	3.517	84	45187	4.516	ng/ul	0.02
7) Phenol-d5	6.775	99	44274	4.114	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	188634	24.721	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	144879	17.158	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	86837	10.277	ng/ul	0.00
21) Nitrobenzene-d5	8.757	128	101610	29.117	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143	81366	25.466	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.010	165	136822	21.205	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	203136	18.008	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	599299	29.846	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	688908	28.502	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	5691m	1.552	ng/ul	0.00
60) Fluorene-d10	15.269	176	494019	28.769	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	49921	18.574	ng/ul	0.00
73) Anthracene-d10	17.133	188	806284	31.001	ng/ul	0.00
81) Pyrene-d10	19.398	212	906013	30.891	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.327	264	818393	30.747	ng/ul	0.00
Target Compounds						Qvalue

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

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