Quantitation Report (Qedit)

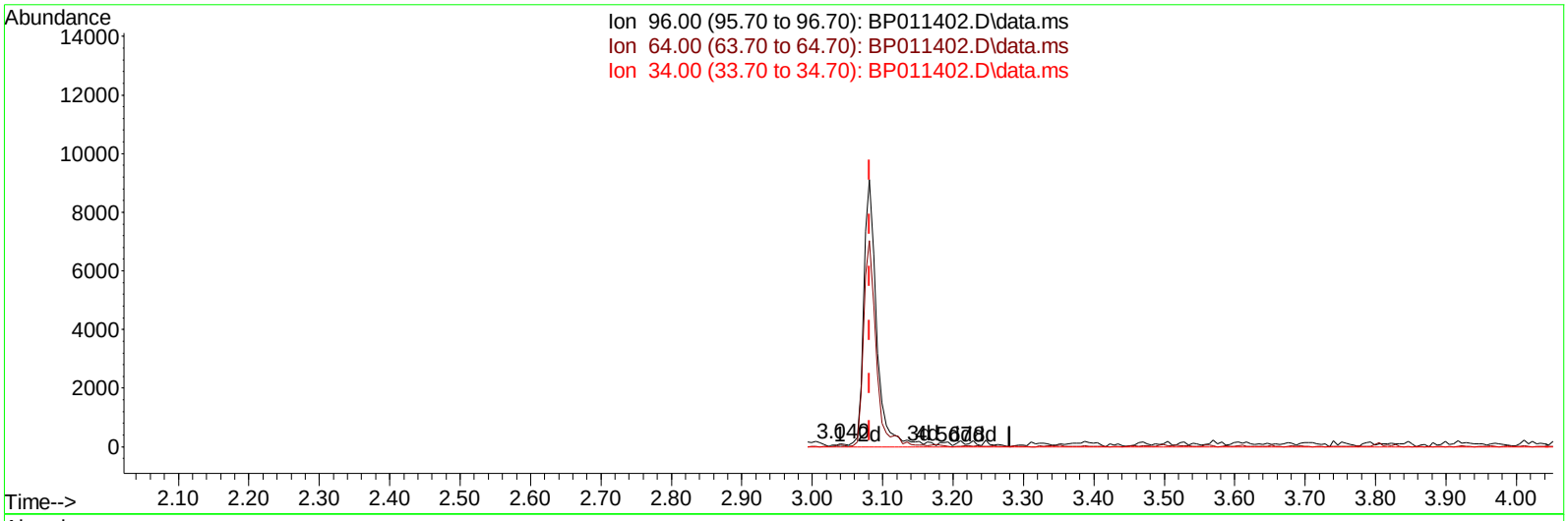
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011402.D
Acq On : 12 Aug 2022 00:16
Operator : CG/JU
Sample : N4128-03
Misc :
ALS Vial : 25 Sample Multiplier: 1

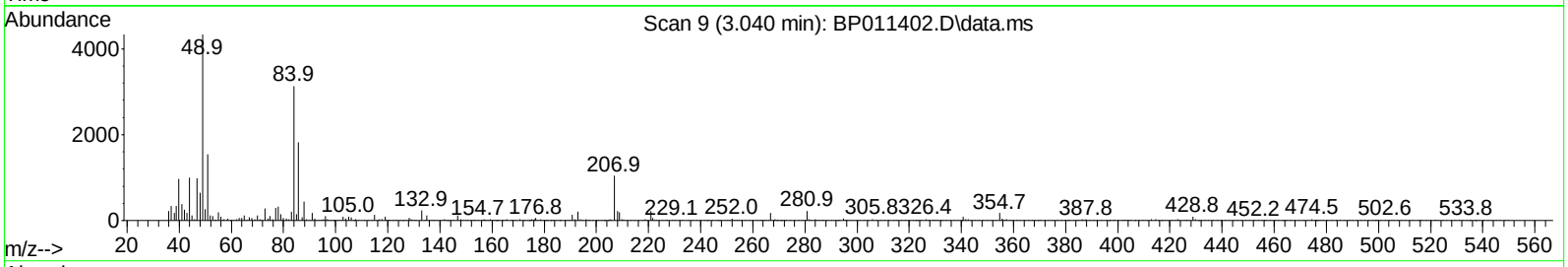
Quant Time: Aug 12 01:11:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 11 23:05:53 2022
Response via : Initial Calibration

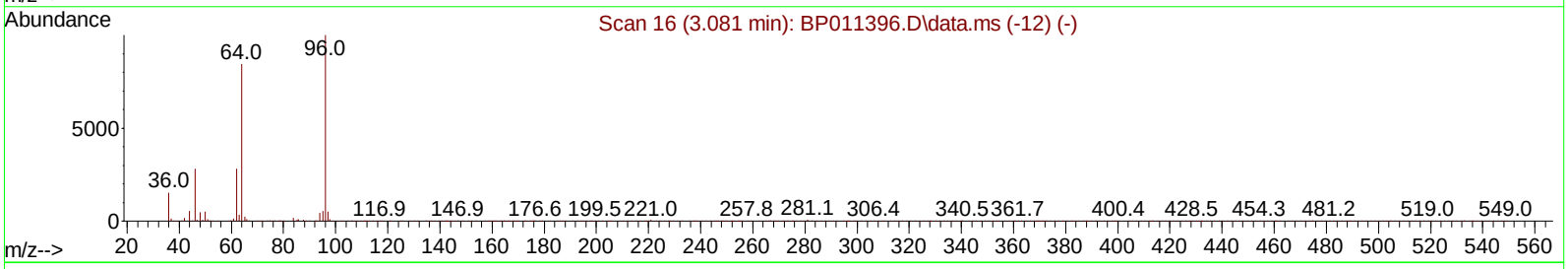
Instrument :
BNA_P
ClientSampleId :
BGBT5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/12/2022
Supervised By :Sohil Jodhani 08/13/2022







TIC: BP011402.D\data.ms

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

3.040min (-0.041) 0.03 ng/uL

response 89

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	49.25
34.00	0.00	0.00
0.00	0.00	0.00

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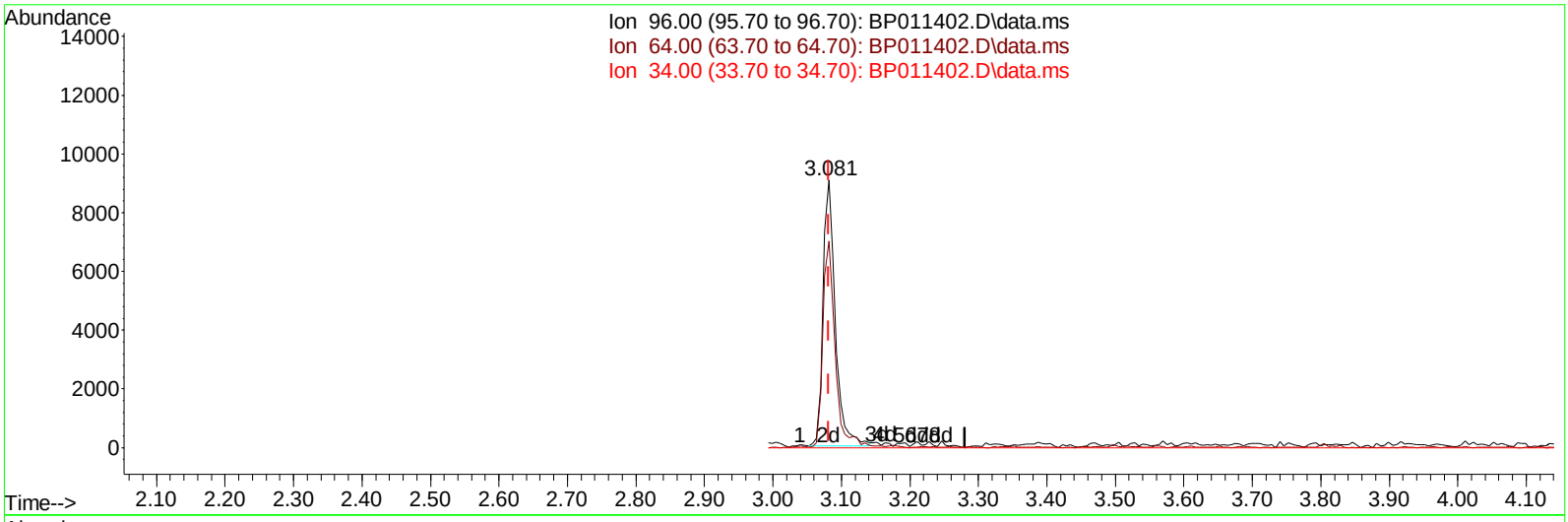

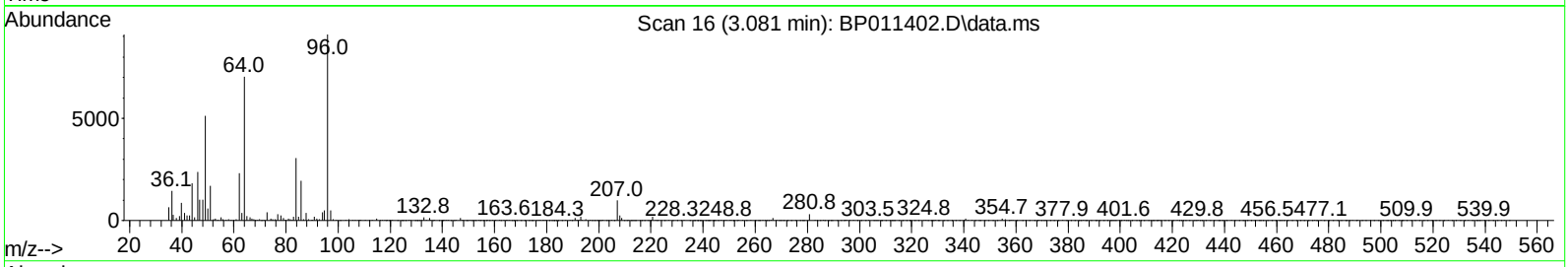

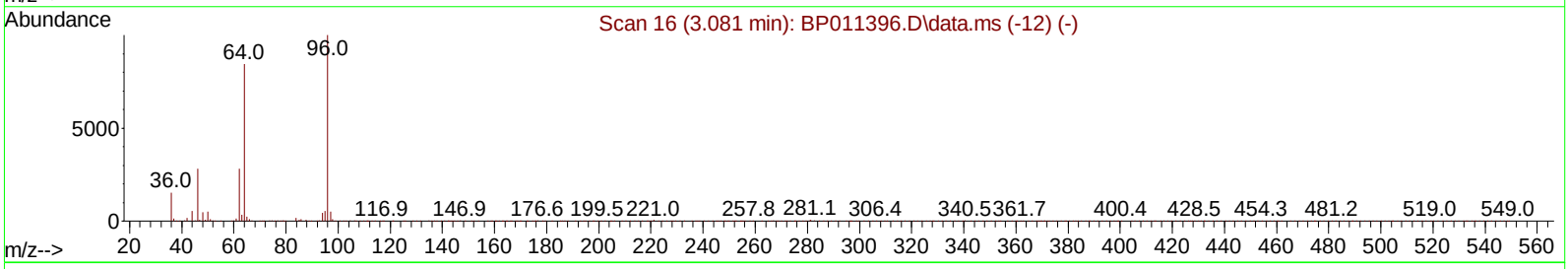

TIC: BP011402.D\data.ms

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

3.081min (+ 0.000) 3.44 ng/uL m

response 11291

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	77.30#
34.00	0.00	0.00
0.00	0.00	0.00

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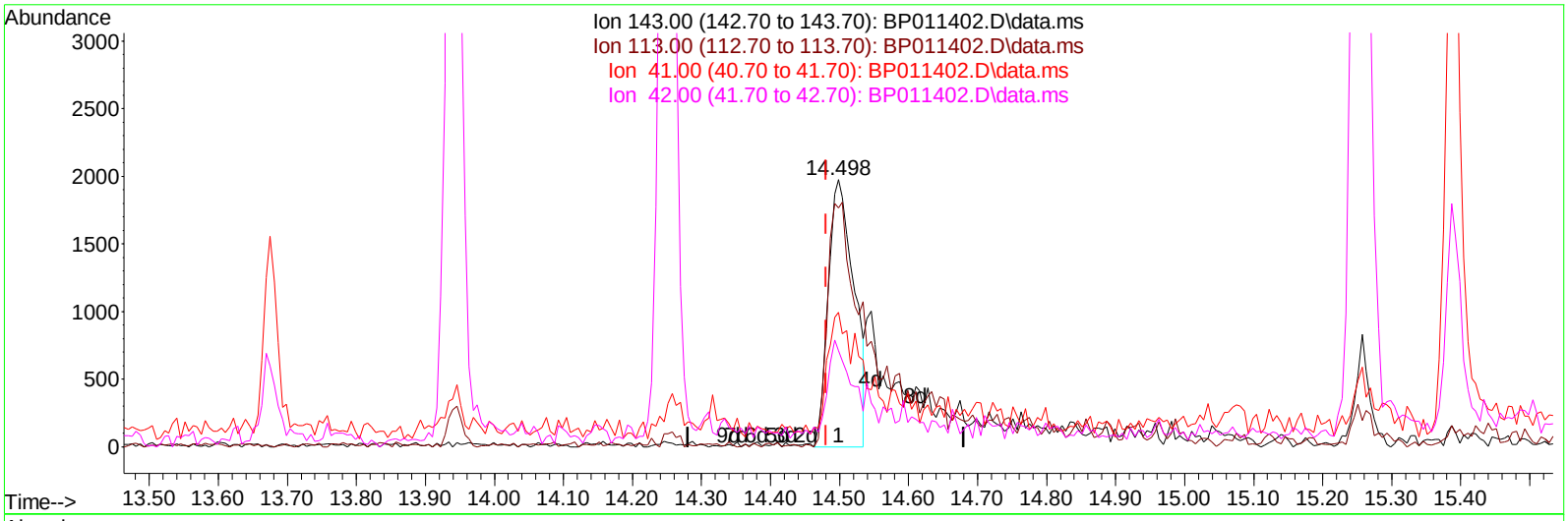

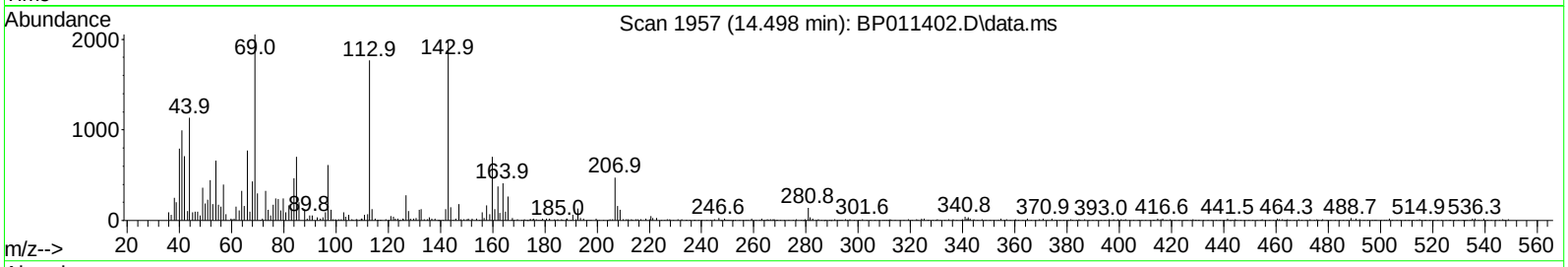

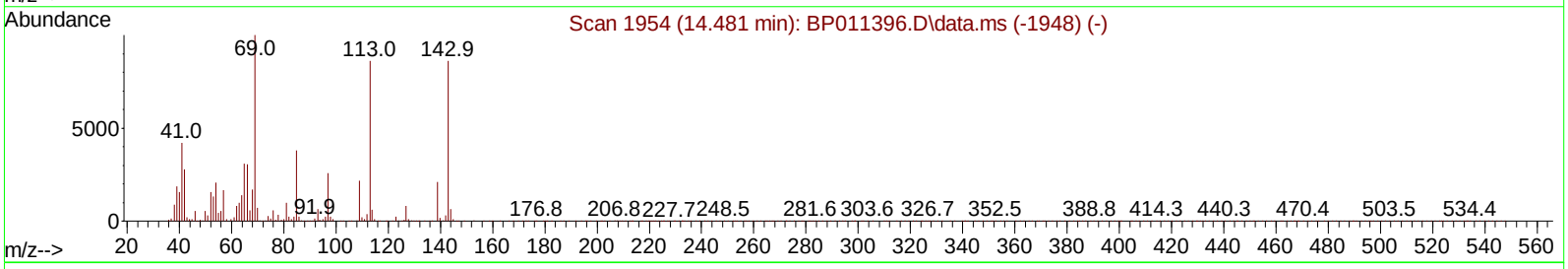

TIC: BP011402.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 1.64 ng/ul

response 5116

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	89.51
41.00	54.40	50.35
42.00	35.50	35.92

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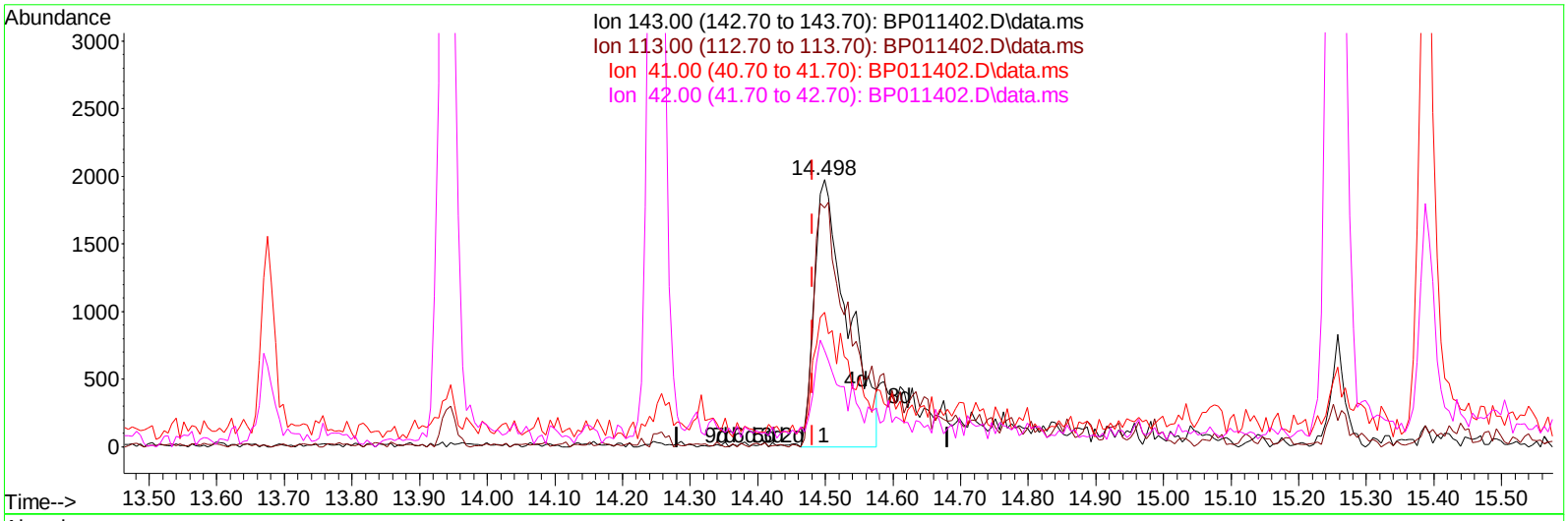

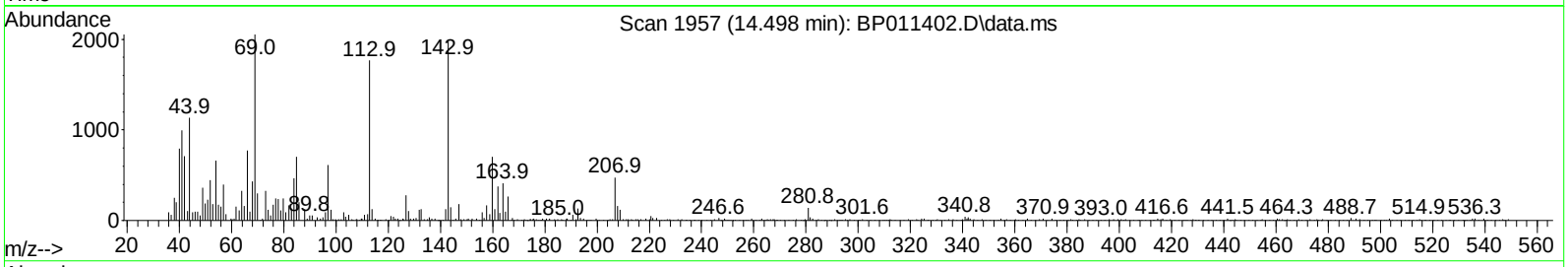

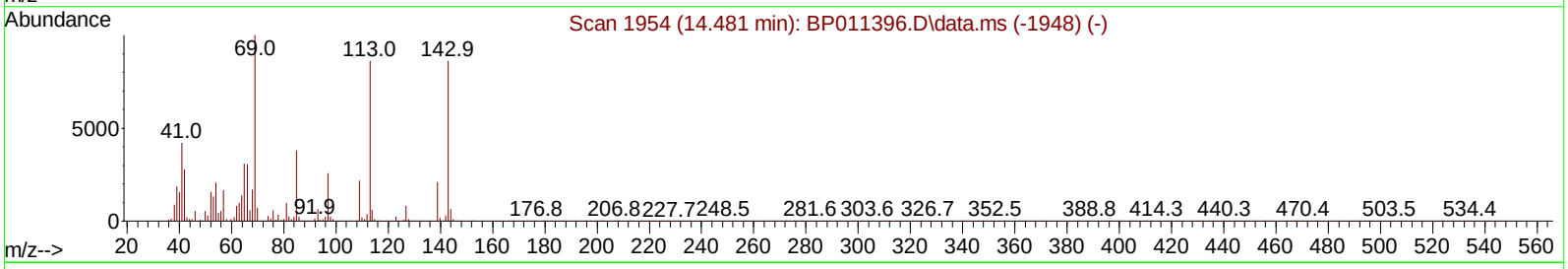

TIC: BP011402.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 2.15 ng/ul m

response 6711

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	89.51
41.00	54.40	50.35
42.00	35.50	35.92

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	100587	20.000	ng/ul	0.00
20) Naphthalene-d8	10.369	136	419198	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.251	164	246224	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	536390	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.157	240	559150	20.000	ng/ul	# 0.00
88) Perylene-d12	23.463	264	609537	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	11291m	3.441	ng/uL	0.00
4) Pyridine-d5	3.499	84	44621	5.391	ng/ul	0.01
7) Phenol-d5	6.764	99	37631	4.227	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	137836	21.838	ng/ul	0.00
11) 2-Chlorophenol-d4	7.111	132	128827	18.445	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	72476	10.369	ng/ul	0.00
21) Nitrobenzene-d5	8.740	128	77538	27.178	ng/ul	0.00
24) 2-Nitrophenol-d4	9.458	143	74046	28.348	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	116346	22.057	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	159802	17.329	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	491522	28.771	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	546781	26.588	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	6711m	2.151	ng/ul	0.02
60) Fluorene-d10	15.257	176	411814	28.186	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.387	200	57923	23.249	ng/ul	0.00
73) Anthracene-d10	17.122	188	737311	30.582	ng/ul	0.00
81) Pyrene-d10	19.386	212	877765	28.494	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.316	264	920979	29.961	ng/ul	0.00
Target Compounds						Qvalue

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

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