Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
Data File : BP010525.D
Acq On : 24 May 2022 20:14
Operator : CG/JU
Sample : N2950-11
Misc :
ALS Vial : 54 Sample Multiplier: 1

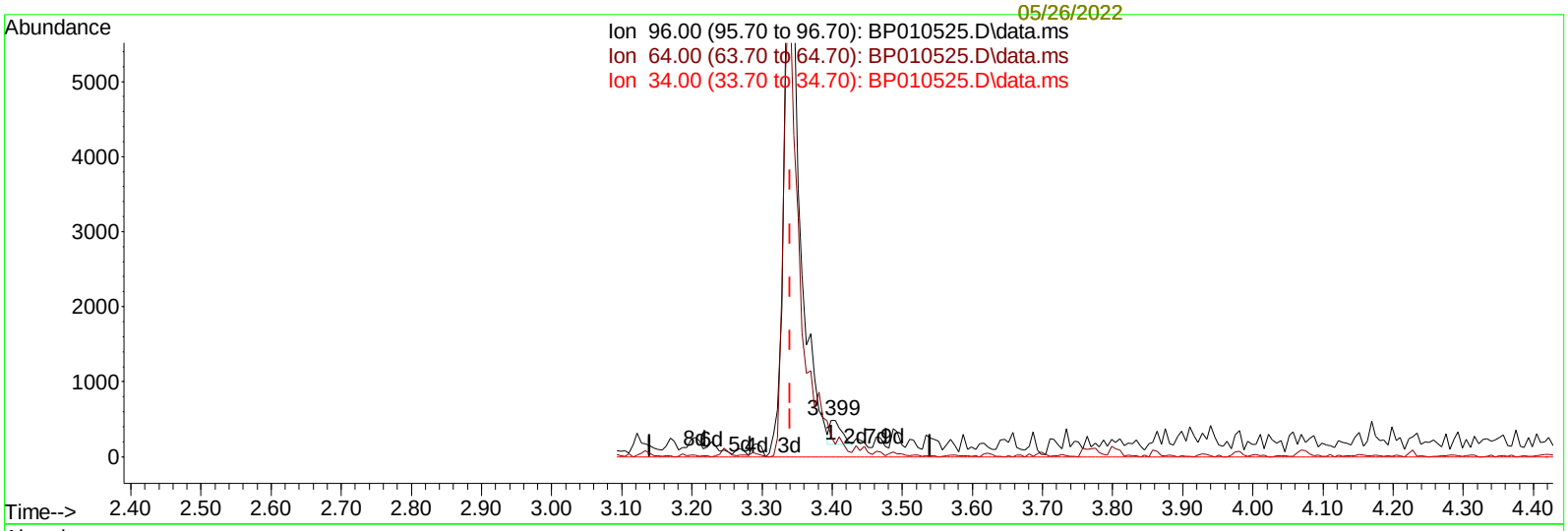
Quant Time: May 24 23:18:22 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 24 05:50:40 2022
Response via : Initial Calibration

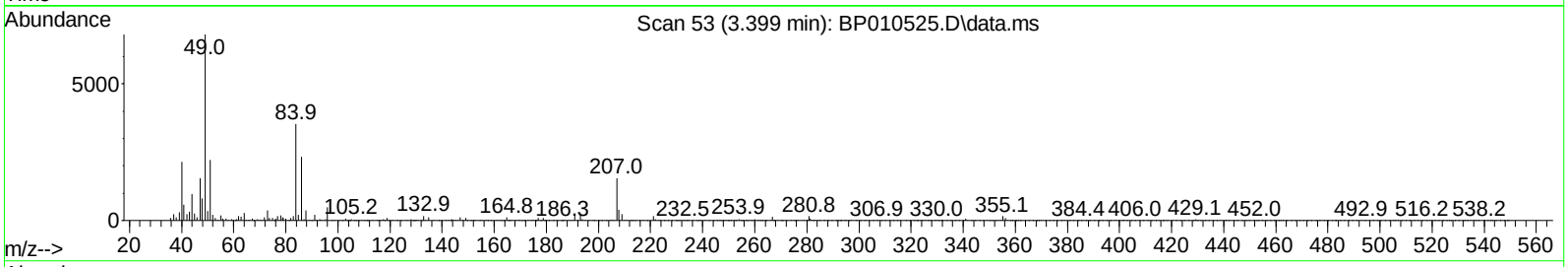
Instrument :
BNA_P
ClientSampleId :
DBTD1

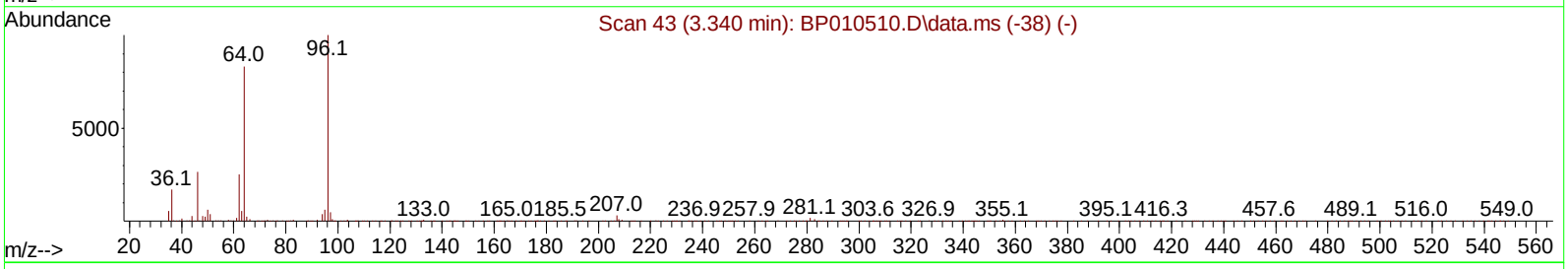
Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay
05/25/2022

Supervised By :Yogesh Patel
05/26/2022







TIC: BP010525.D\data.ms

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

3.399min (+ 0.058) 0.13 ng/uL

response 363

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.17
34.00	0.00	0.00
0.00	0.00	0.00

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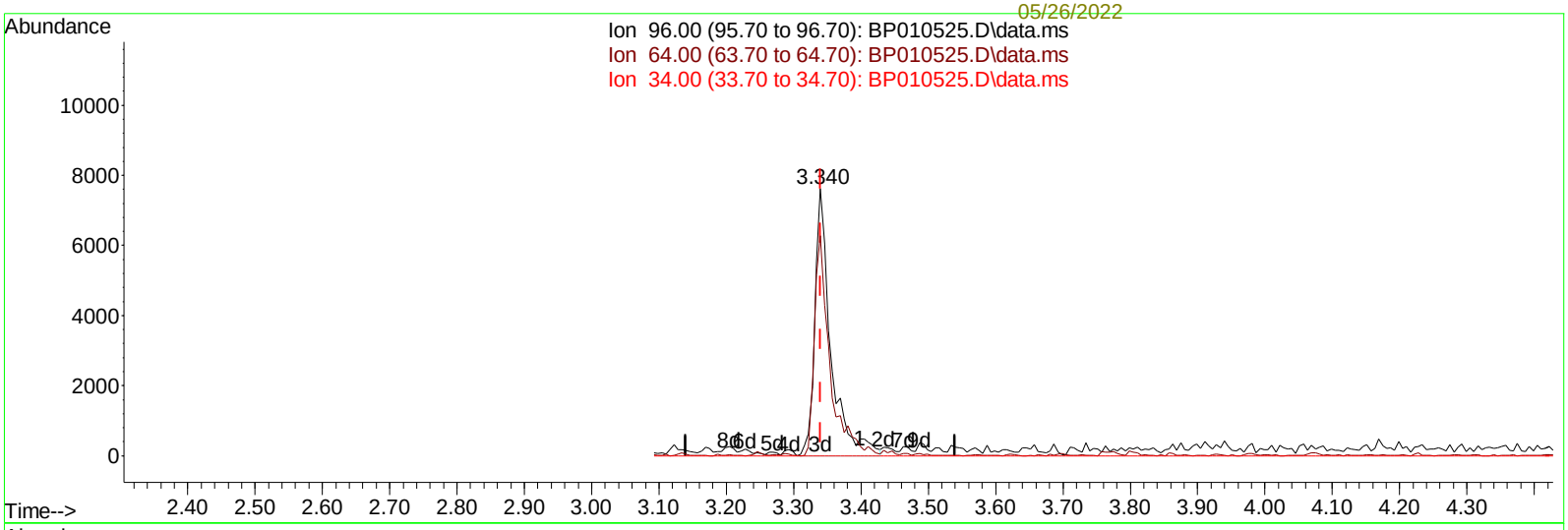

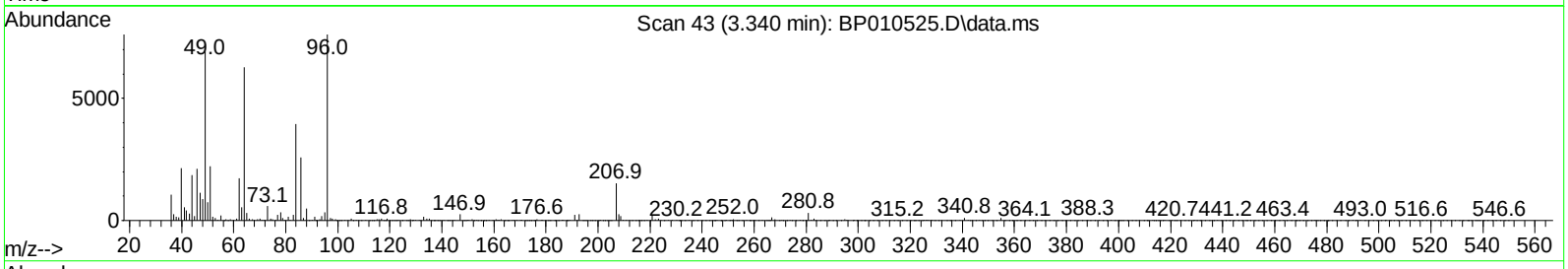

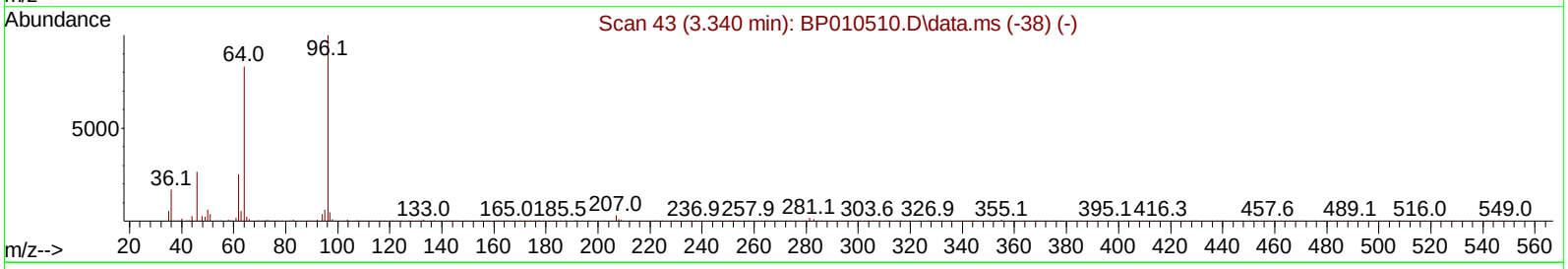

TIC: BP010525.D\data.ms

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

3.340min (-0.000) 4.59 ng/uL m

response 12632

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	82.46#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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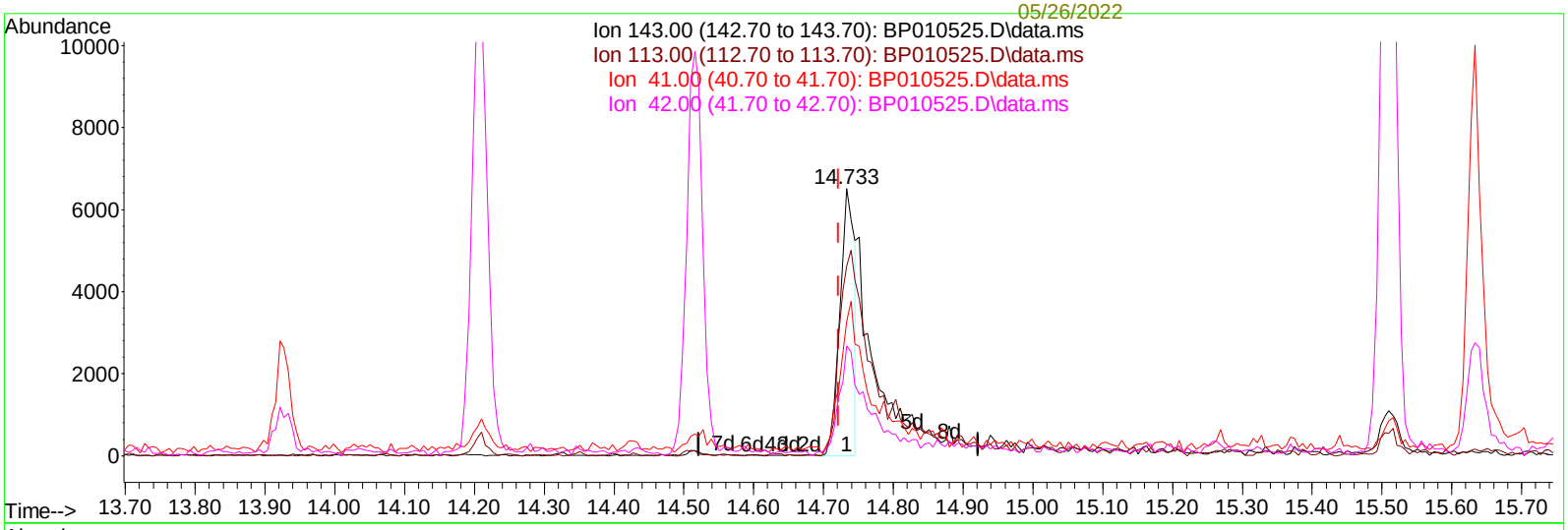

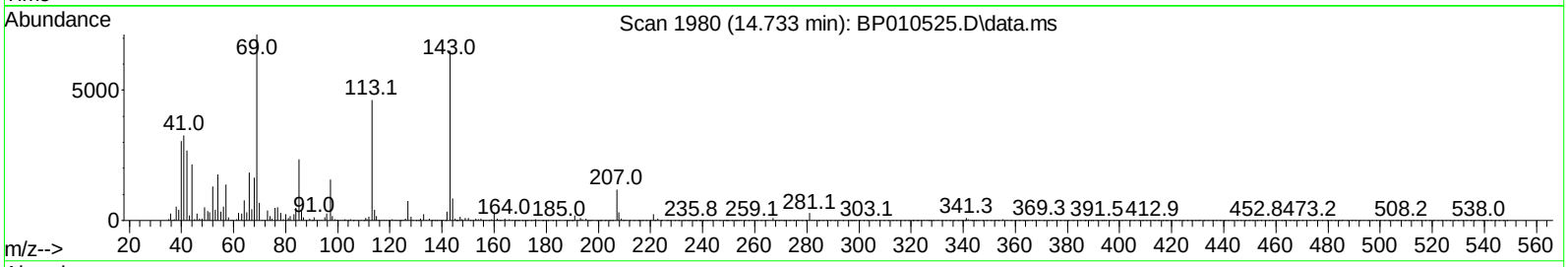

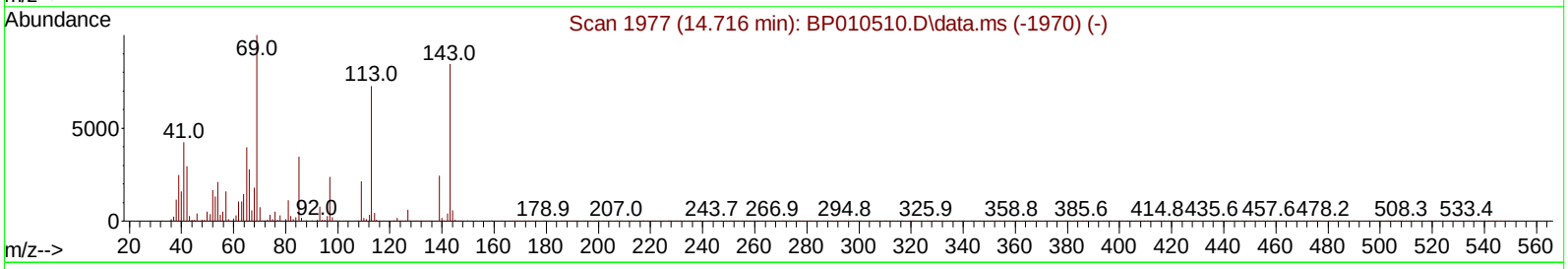

TIC: BP010525.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.011) 1.85 ng/ul

response 9547

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	71.04#
41.00	23.20	50.30#
42.00	18.00	41.14#

Quantitation Report (Qedit)

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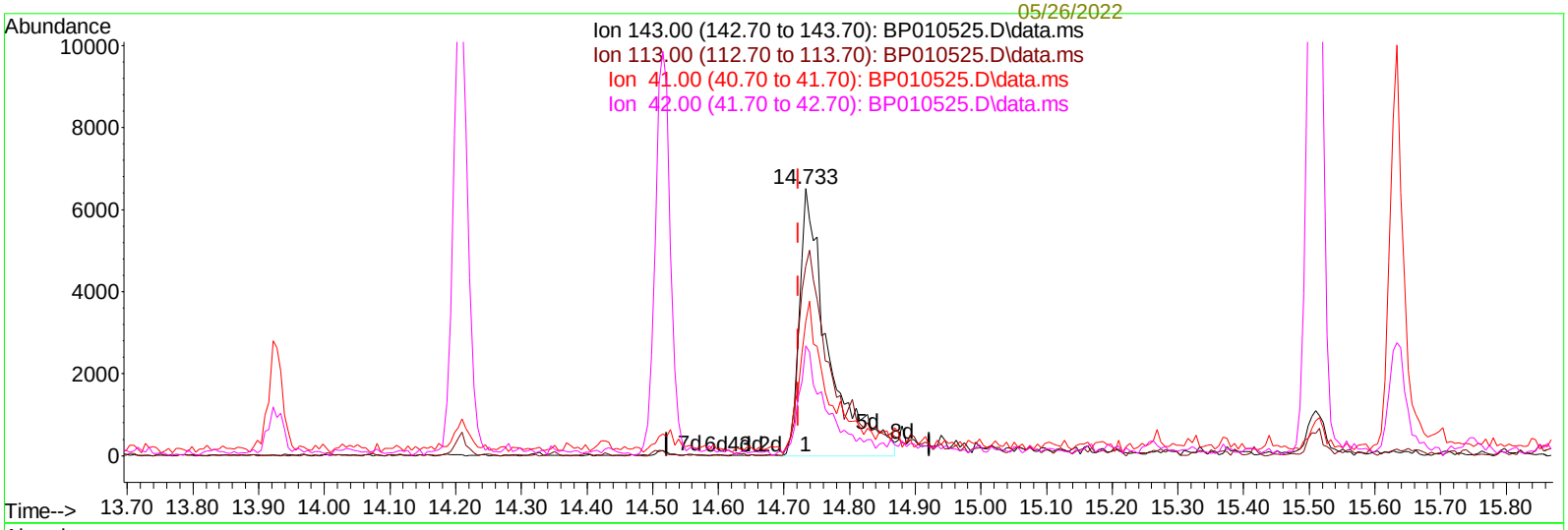

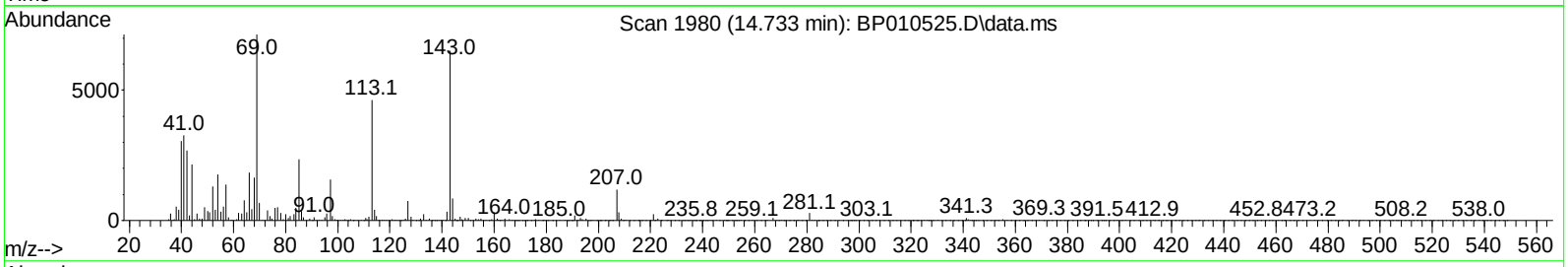

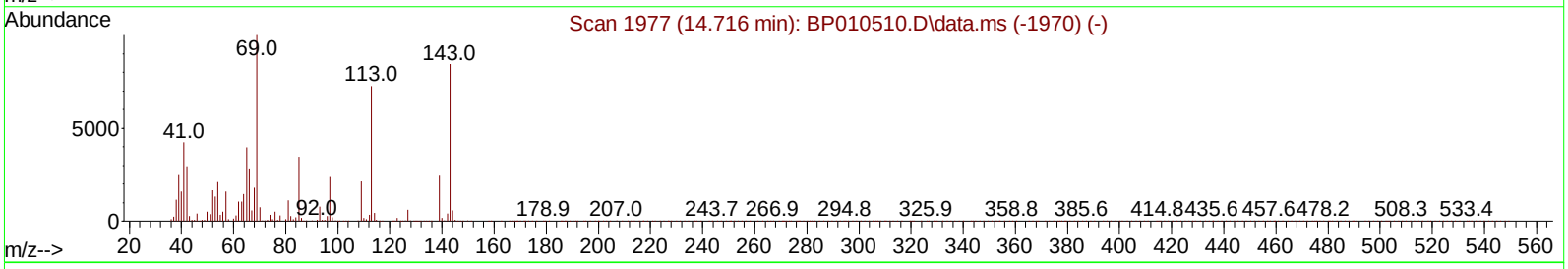

TIC: BP010525.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.011) 3.90 ng/ul m

response 20080

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	71.04#
41.00	23.20	50.30#
42.00	18.00	41.14#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	117179	20.000	ng/ul	0.00
20) Naphthalene-d8	10.686	136	559921	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	393019	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.268	188	860573	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.351	240	770870	20.000	ng/ul	# 0.00
88) Perylene-d12	23.774	264	622296	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	12632m	4.595	ng/uL	0.00
4) Pyridine-d5	3.769	84	56303	7.213	ng/ul	0.01
7) Phenol-d5	7.051	99	63138	6.420	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	220005	33.294	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	190796	25.352	ng/ul	0.00
15) 4-Methylphenol-d8	8.592	113	133137	15.929	ng/ul	0.00
21) Nitrobenzene-d5	9.045	128	143515	33.118	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	146703	31.945	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	242412	27.916	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	322125	24.949	ng/ul	0.00
46) Dimethylphthalate-d6	13.927	166	980119	33.848	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	1115506	33.398	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	20080m	3.896	ng/ul	0.01
60) Fluorene-d10	15.510	176	869603	34.520	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	124742	24.785	ng/ul	0.00
73) Anthracene-d10	17.368	188	1504622	38.612	ng/ul	0.00
81) Pyrene-d10	19.598	212	1807158	44.143	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.615	264	1241972	39.746	ng/ul	0.00

Target Compounds Qvalue

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

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