Quantitation Report (Qedit)

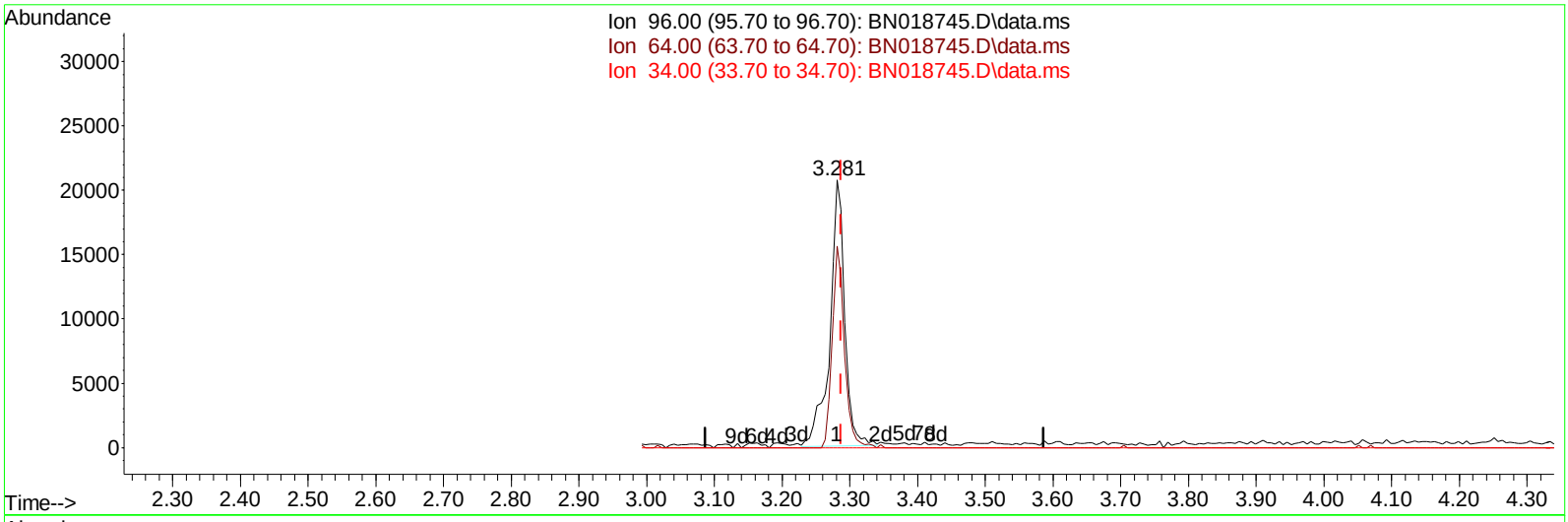
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022522\
Data File : BN018745.D
Acq On : 26 Feb 2022 00:15
Operator : CG/JU
Sample : N1656-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

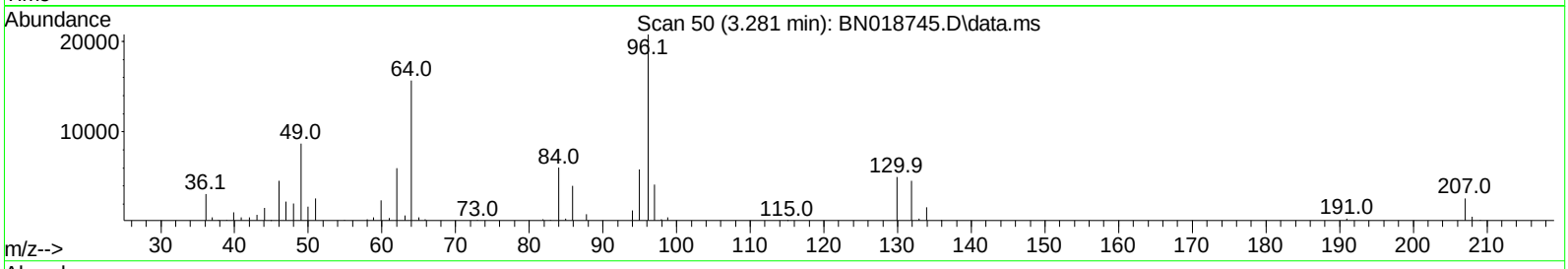
Quant Time: Feb 26 01:14:09 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Feb 16 16:48:14 2022
Response via : Initial Calibration

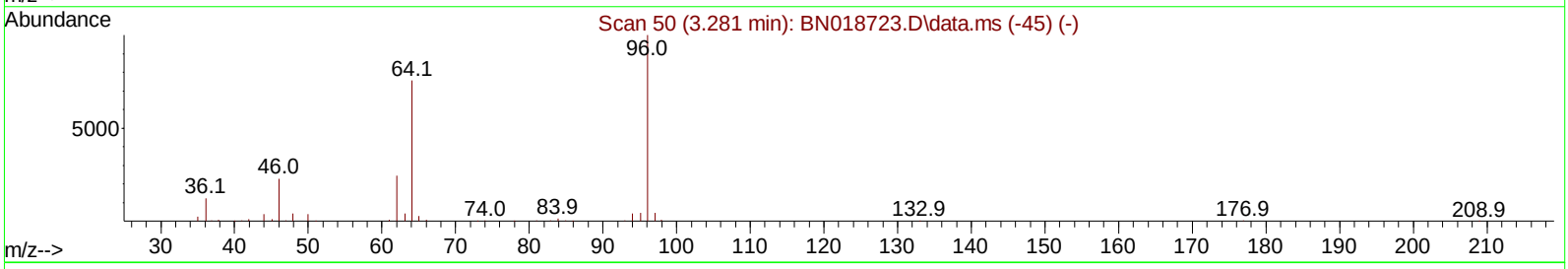
Instrument :
BNA_N
ClientSampleId :
BGQ23

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/28/2022
Supervised By :mohammad ahmed 02/28/2022







TIC: BN018745.D\data.ms

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

3.281min (-0.006) 5.85 ng/uL

response 32050

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.80	75.30
34.00	0.00	0.00
0.00	0.00	0.00

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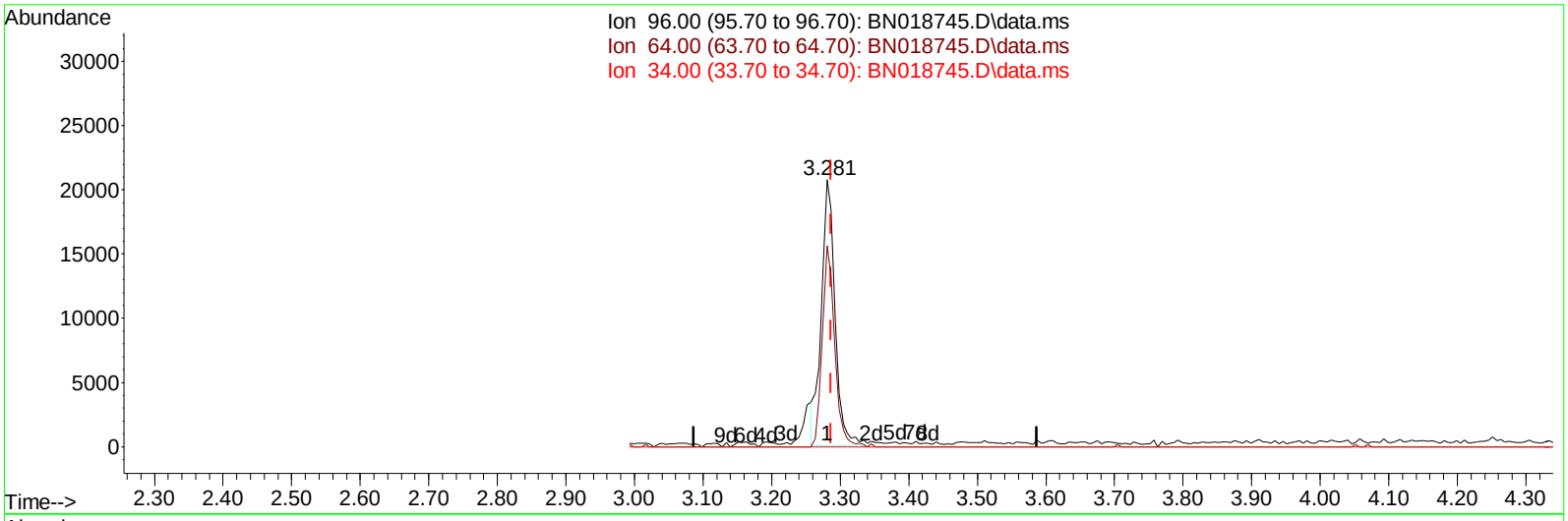

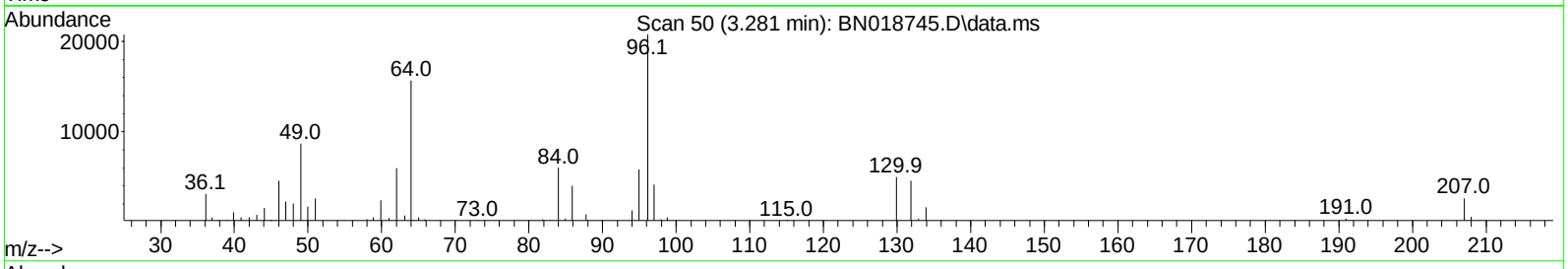

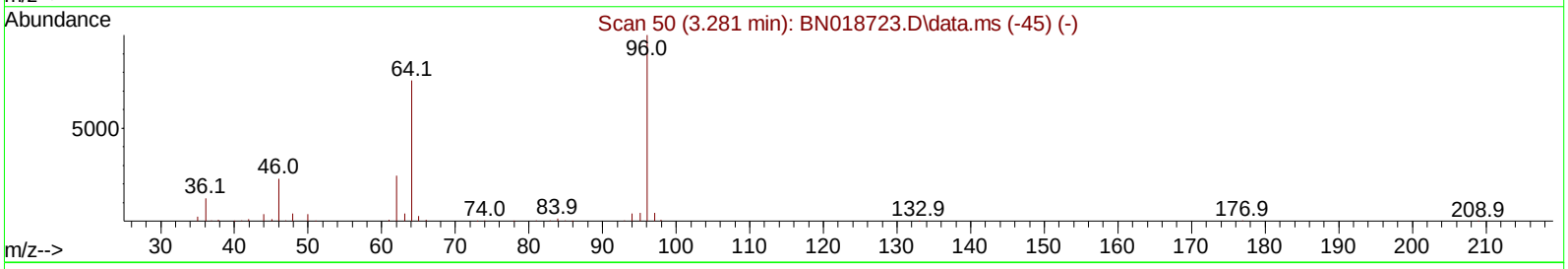

TIC: BN018745.D\data.ms

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

3.281min (-0.006) 5.29 ng/uL m

response 28963

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.80	75.30
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

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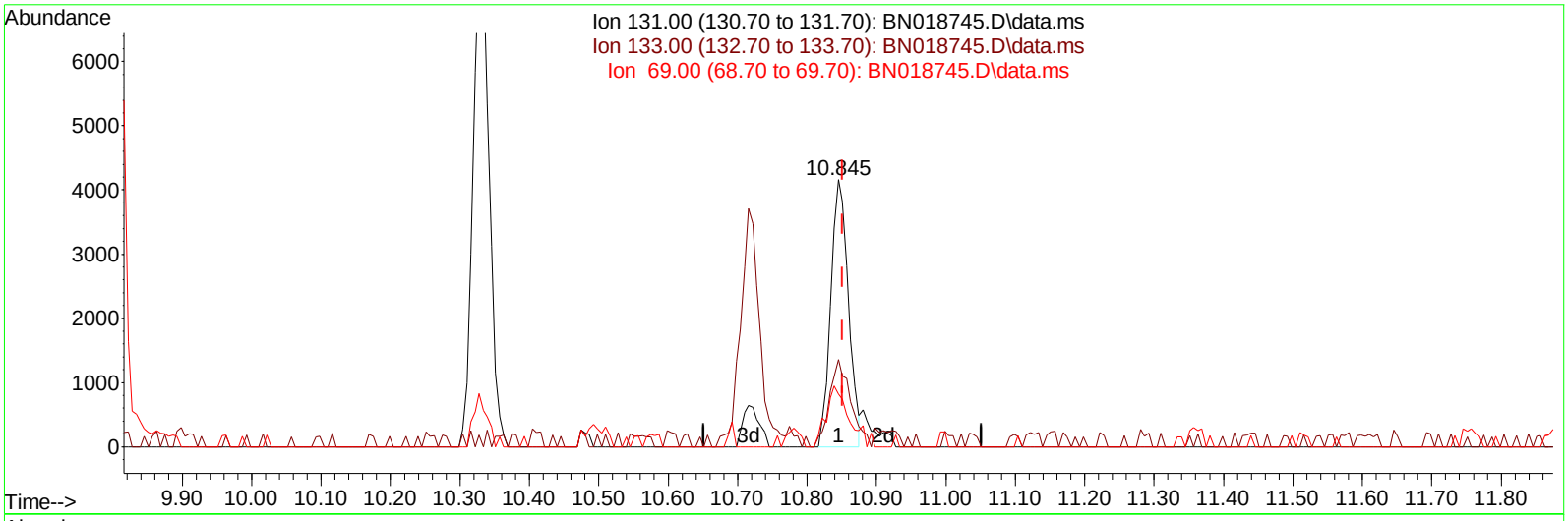

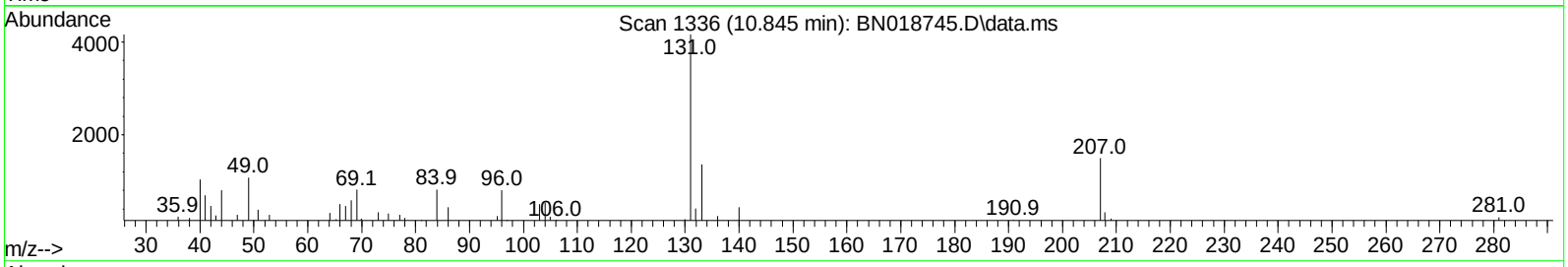

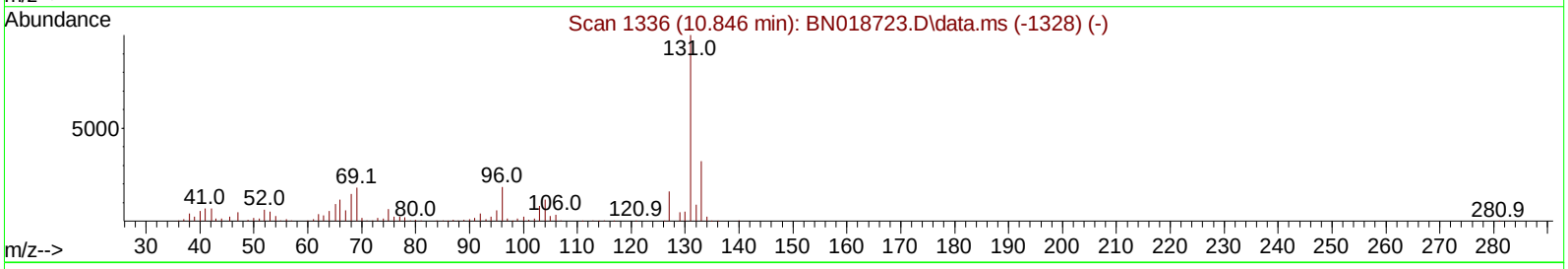

TIC: BN018745.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.845min (-0.006) 0.45 ng/ul

response 7370

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.30	32.73
69.00	20.30	19.67
0.00	0.00	0.00

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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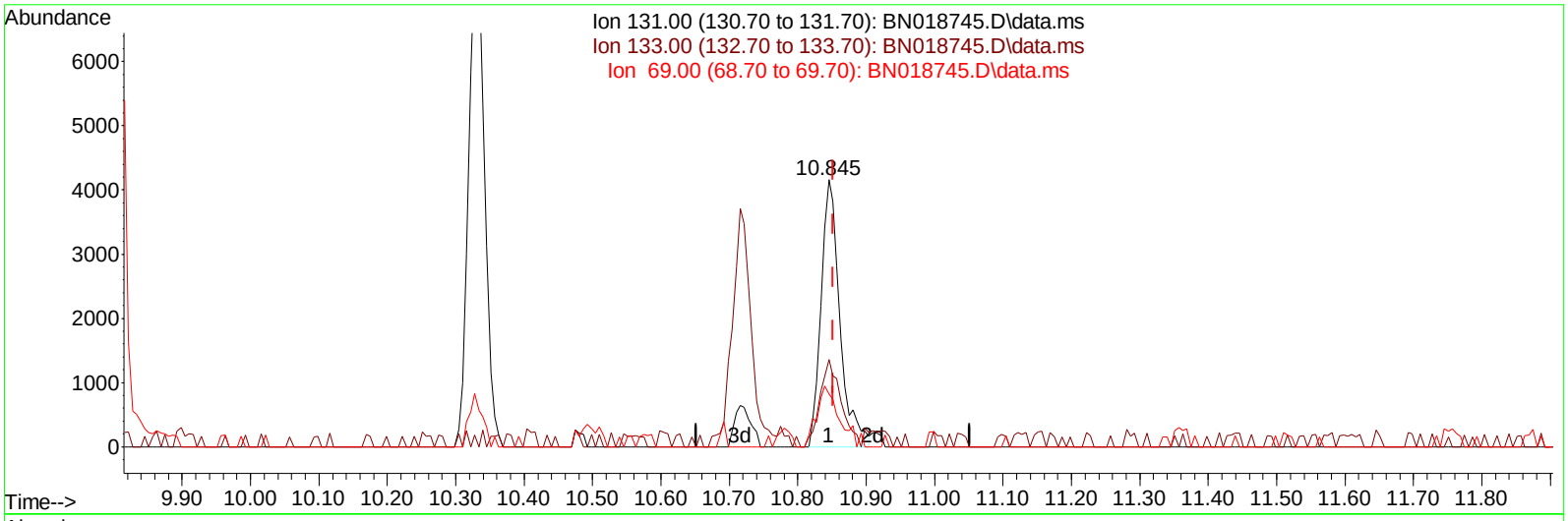

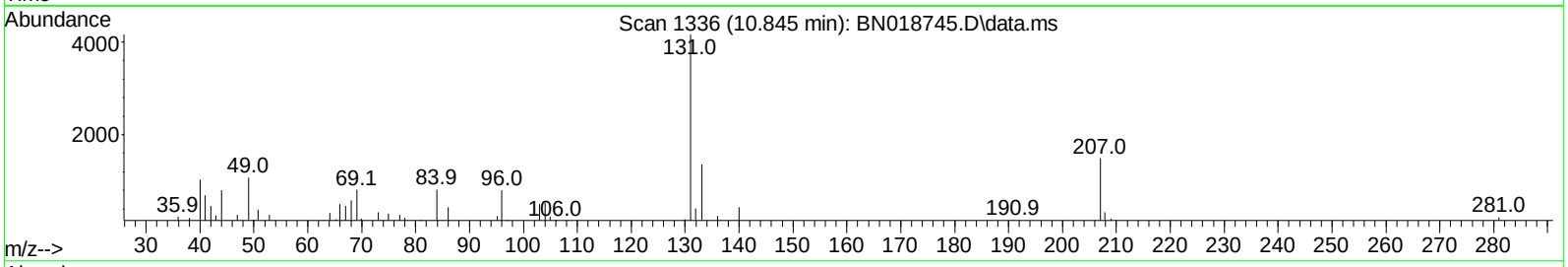

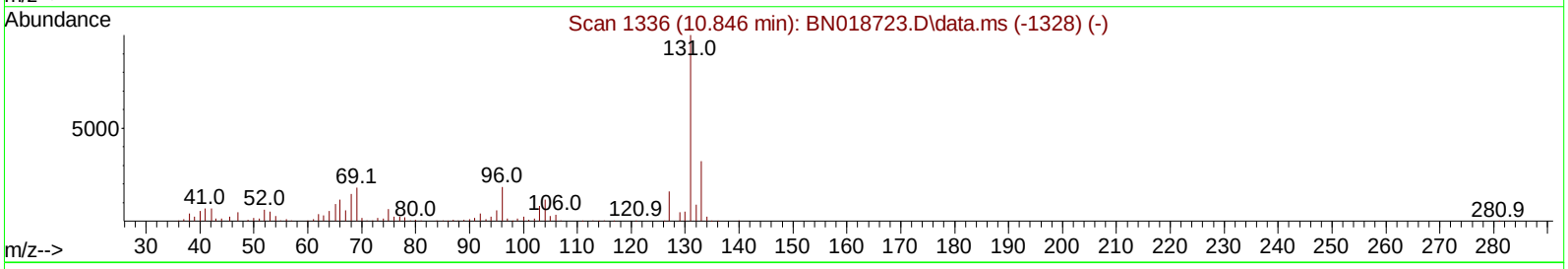

TIC: BN018745.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.845min (-0.006) 0.49 ng/ul m

response 7941

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.30	32.73
69.00	20.30	19.67
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022522\
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/28/2022
Supervised By :mohammad ahmed 02/28/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	206338	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	776331	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.545	164	440588	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	831132	20.000	ng/ul	0.00
79) Chrysene-d12	21.462	240	754116	20.000	ng/ul	0.00
88) Perylene-d12	23.856	264	776525	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	28963m	5.291	ng/uL	0.00
4) Pyridine-d5	3.710	84	35896	2.425	ng/ul	0.00
7) Phenol-d5	7.057	99	133018	7.816	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.228	67	335203	31.737	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	350062	26.118	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	223977	17.235	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	191838	32.158	ng/ul	0.00
24) 2-Nitrophenol-d4	9.792	143	201110	31.478	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	350387	28.734	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	7941m	0.487	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1151545	36.104	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	1413412	34.205	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	33988	6.132	ng/ul	0.00
60) Fluorene-d10	15.533	176	997887	36.239	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	143058	29.173	ng/ul	0.00
73) Anthracene-d10	17.380	188	1590056	39.982	ng/ul	0.00
81) Pyrene-d10	19.674	212	1723545	39.048	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.703	264	1574909	39.769	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.316	88	14212	2.457	ng/uL#	92

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

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