

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092320\
 Data File : BP003365.D
 Acq On : 23 Sep 2020 19:30
 Operator : CG/JU
 Sample : L4053-05DL2 100X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 1799-1801DL2

Quant Time: Sep 24 09:50:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 12:48:08 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	191769	20.00	ng	0.00
21) Naphthalene-d8	10.345	136	780566	20.00	ng	# 0.00
39) Acenaphthene-d10	14.216	164	472301	20.00	ng	0.00
64) Phenanthrene-d10	16.969	188	964702	20.00	ng	0.00
76) Chrysene-d12	21.068	240	773816	20.00	ng	0.00
86) Perylene-d12	23.257	264	713881	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	6762	0.62	ng	0.00
7) Phenol-d6	6.775	99	5749	0.39	ng	0.00
23) Nitrobenzene-d5	8.716	82	11981	0.90	ng	0.00
42) 2,4,6-Tribromophenol	15.722	330	7240	1.01	ng	0.00
45) 2-Fluorobiphenyl	12.834	172	31167	0.97	ng	0.00
79) Terphenyl-d14	19.551	244	94961	2.56	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	8.351	107	743354	55.25	ng	Qvalue 95
32) Benzoic acid	9.751	122	191087	30.20	ng	84

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

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