

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070820\
 Data File : BP002683.D
 Acq On : 08 Jul 2020 19:43
 Operator : CG/JU
 Sample : L3203-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RO-06

Quant Time: Jul 09 08:00:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	204850	20.00	ng	-0.01
21) Naphthalene-d8	10.34	136	865394	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	551707	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	1176229	20.00	ng	0.00
76) Chrysene-d12	21.09	240	1389132	20.00	ng	0.00
86) Perylene-d12	23.28	264	1292922	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	1085337	89.70	ng	0.00
7) Phenol-d6	6.77	99	1629335	96.48	ng	0.00
23) Nitrobenzene-d5	8.71	82	1028461	61.23	ng	-0.01
42) 2,4,6-Tribromophenol	15.72	330	561436	83.67	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	2210741	58.98	ng	0.00
79) Terphenyl-d14	19.57	244	3411108	53.00	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.68	163	319585	7.375	ng	99
70) Pentachlorophenol	16.65	266	291467	37.480	ng	98
71) Phenanthrene	17.02	178	150844	2.384	ng	98
75) Fluoranthene	19.00	202	171405	2.238	ng	96

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

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