

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070620\
 Data File : BN011102.D
 Acq On : 06 Jul 2020 11:48
 Operator : CG/JU
 Sample : PB129903BL
 Misc : MDL-WTR-PBBL-04
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB129903BL

Quant Time: Jul 06 12:32:27 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN063020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 06 10:52:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	1935	0.40	ng	0.00
7) Naphthalene-d8	10.28	136	5823	0.40	ng	0.01
13) Acenaphthene-d10	14.14	164	3052	0.40	ng	0.01
19) Phenanthrene-d10	16.91	188	6331	0.40	ng	0.02
29) Chrysene-d12	21.11	240	5268	0.40	ng	0.01
36) Perylene-d12	23.25	264	5876	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	1285	0.27	ng	0.00
5) Phenol-d6	6.74	99	1327	0.24	ng	0.02
8) Nitrobenzene-d5	8.68	82	1452	0.29	ng	0.03
11) 2-Methylnaphthalene-d10	11.88	152	3591	0.36	ng	0.02
14) 2,4,6-Tribromophenol	15.68	330	308	0.24	ng	0.04
15) 2-Fluorobiphenyl	12.77	172	3799	0.26	ng	0.03
27) Fluoranthene-d10	18.94	212	7624	0.40	ng	0.00
31) Terphenyl-d14	19.55	244	5762	0.42	ng	0.00

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

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