

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070720\
 Data File : BN011115.D
 Acq On : 07 Jul 2020 14:44
 Operator : CG/JU
 Sample : PB129932BL
 Misc : MDL-WTR-PBBL-07
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB129932BL

Quant Time: Jul 07 15:16:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN063020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 07 14:39:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	2049	0.40	ng	0.00
7) Naphthalene-d8	10.28	136	5691	0.40	ng	0.03
13) Acenaphthene-d10	14.14	164	2905	0.40	ng	0.01
19) Phenanthrene-d10	16.91	188	5983	0.40	ng	0.02
29) Chrysene-d12	21.11	240	5185	0.40	ng	0.01
36) Perylene-d12	23.24	264	5400	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	1273	0.25	ng	0.00
5) Phenol-d6	6.73	99	1139	0.19	ng	0.00
8) Nitrobenzene-d5	8.68	82	1385	0.29	ng	0.03
11) 2-Methylnaphthalene-d10	11.89	152	3467	0.35	ng	0.03
14) 2,4,6-Tribromophenol	15.68	330	299	0.25	ng	0.04
15) 2-Fluorobiphenyl	12.77	172	3604	0.26	ng	0.03
27) Fluoranthene-d10	18.94	212	7360	0.41	ng	0.00
31) Terphenyl-d14	19.55	244	5252	0.39	ng	0.00

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

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