

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031122\
 Data File : BN018906.D
 Acq On : 10 Mar 2022 17:14
 Operator : CG/JU
 Sample : PB143174BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143174BL

Quant Time: Mar 11 03:04:54 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN030322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 03:02:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.912	152	4588	0.400	ng	0.00
7) Naphthalene-d8	10.712	136	12765	0.400	ng	# 0.00
13) Acenaphthene-d10	14.538	164	8028	0.400	ng	0.00
19) Phenanthrene-d10	17.277	188	19140	0.400	ng	0.00
29) Chrysene-d12	21.458	240	17910	0.400	ng	0.00
35) Perylene-d12	23.849	264	15479	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.464	112	3275	0.326	ng	0.00
5) Phenol-d6	7.067	99	3661	0.304	ng	0.00
8) Nitrobenzene-d5	9.067	82	2525	0.271	ng	0.01
11) 2-Methylnaphthalene-d10	12.302	152	5655	0.272	ng	0.00
14) 2,4,6-Tribromophenol	16.025	330	703	0.190	ng	0.00
15) 2-Fluorobiphenyl	13.169	172	9362	0.272	ng	0.00
27) Fluoranthene-d10	19.303	212	15769	0.280	ng	0.00
31) Terphenyl-d14	19.898	244	11215	0.293	ng	0.00

Target Compounds Qvalue

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

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