

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032122\
 Data File : BN019035.D
 Acq On : 21 Mar 2022 23:55
 Operator : CG/JU
 Sample : PB143429BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143429BL

Quant Time: Mar 22 02:33:23 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 13:18:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	6715	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	17248	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	8720	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	11425	0.400	ng	0.00
29) Chrysene-d12	21.422	240	8564	0.400	ng	# 0.00
35) Perylene-d12	23.787	264	9014	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	5699	0.379	ng	0.00
5) Phenol-d6	7.017	99	4671	0.296	ng	0.00
8) Nitrobenzene-d5	9.014	82	3925	0.316	ng	0.00
11) 2-Methylnaphthalene-d10	12.250	152	8562	0.299	ng	0.00
14) 2,4,6-Tribromophenol	15.995	330	575	0.137	ng	0.01
15) 2-Fluorobiphenyl	13.127	172	11559	0.331	ng	0.00
27) Fluoranthene-d10	19.265	212	10970	0.324	ng	0.00
31) Terphenyl-d14	19.860	244	6227	0.296	ng	0.00

Target Compounds Qvalue

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

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