

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032122\
 Data File : BN019019.D
 Acq On : 21 Mar 2022 12:22
 Operator : CG/JU
 Sample : PB143448BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143448BL

Quant Time: Mar 22 02:29:17 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	7.847	152	6718	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	16976	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	9317	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	14897	0.400	ng	0.00
29) Chrysene-d12	21.431	240	10047	0.400	ng	# 0.00
35) Perylene-d12	23.799	264	7422	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.399	112	5066	0.337	ng	0.00
5) Phenol-d6	7.009	99	4045	0.256	ng	0.00
8) Nitrobenzene-d5	9.014	82	3456	0.282	ng	0.00
11) 2-Methylnaphthalene-d10	12.250	152	8669	0.308	ng	0.00
14) 2,4,6-Tribromophenol	15.995	330	549	0.122	ng	0.01
15) 2-Fluorobiphenyl	13.116	172	12212	0.327	ng	-0.01
27) Fluoranthene-d10	19.269	212	15845	0.358	ng	0.00
31) Terphenyl-d14	19.864	244	9503	0.385	ng	0.00

Target Compounds Qvalue

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

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