

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092521\
 Data File : BN016468.D
 Acq On : 25 Sep 2021 10:53
 Operator : CG/JU
 Sample : PB139340BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB139340BL

Quant Time: Sep 26 04:14:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 03:35:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.670	152	21079	0.40	ng	0.00
7) Naphthalene-d8	10.432	136	89180	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	42792	0.40	ng	0.00
19) Phenanthrene-d10	17.042	188	78476	0.40	ng	0.00
29) Chrysene-d12	21.249	240	59358	0.40	ng	# 0.00
35) Perylene-d12	23.519	264	47794	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.309	112	17454	0.37	ng	0.03
5) Phenol-d6	6.859	99	18405	0.33	ng	0.00
8) Nitrobenzene-d5	8.810	82	24490	0.39	ng	0.00
11) 2-Methylnaphthalene-d10	12.029	152	52411	0.40	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	2652	0.29	ng	0.00
15) 2-Fluorobiphenyl	12.911	172	67470	0.38	ng	-0.01
27) Fluoranthene-d10	19.088	212	79619	0.37	ng	0.00
31) Terphenyl-d14	19.698	244	63147	0.46	ng	0.00

Target Compounds Qvalue

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

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