

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110921\
 Data File : BN017361.D
 Acq On : 09 Nov 2021 23:23
 Operator : CG/JU
 Sample : PB140411BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB140411BL

Quant Time: Nov 10 11:21:53 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 10 11:00:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	31086	0.400	ng	0.00
7) Naphthalene-d8	10.319	136	132256	0.400	ng	0.01
13) Acenaphthene-d10	14.181	164	73151	0.400	ng	0.00
19) Phenanthrene-d10	16.934	188	132675	0.400	ng	0.00
29) Chrysene-d12	21.144	240	114023	0.400	ng	0.01
35) Perylene-d12	23.335	264	121905	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.191	112	23677	0.330	ng	0.03
5) Phenol-d6	6.749	99	21020	0.266	ng	0.02
8) Nitrobenzene-d5	8.697	82	22612	0.285	ng	0.01
11) 2-Methylnaphthalene-d10	11.911	152	69581	0.357	ng	0.00
14) 2,4,6-Tribromophenol	15.691	330	3359	0.123	ng	0.01
15) 2-Fluorobiphenyl	12.811	172	89956	0.323	ng	0.01
27) Fluoranthene-d10	18.975	212	116281	0.344	ng	0.00
31) Terphenyl-d14	19.590	244	97584	0.384	ng	0.00

Target Compounds Qvalue

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

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