

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100422\
 Data File : BN022011.D
 Acq On : 04 Oct 2022 18:26
 Operator : CG/JU
 Sample : PB148092BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB148092BL

Quant Time: Oct 04 23:06:04 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 04 01:32:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.986	152	4336	0.400	ng	0.00
7) Naphthalene-d8	10.816	136	12639	0.400	ng	0.00
13) Acenaphthene-d10	14.653	164	6317	0.400	ng	0.00
19) Phenanthrene-d10	17.412	188	12591	0.400	ng	# 0.00
29) Chrysene-d12	21.609	240	9051	0.400	ng	0.00
35) Perylene-d12	24.087	264	6965	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.516	112	4207	0.419	ng	0.00
5) Phenol-d6	7.148	99	4873	0.374	ng	0.00
8) Nitrobenzene-d5	9.172	82	3566	0.351	ng	0.01
11) 2-Methylnaphthalene-d10	12.414	152	8521	0.367	ng	0.00
14) 2,4,6-Tribromophenol	16.158	330	645	0.268	ng	0.01
15) 2-Fluorobiphenyl	13.285	172	10469	0.379	ng	0.00
27) Fluoranthene-d10	19.445	212	12088	0.361	ng	0.00
31) Terphenyl-d14	20.042	244	7874	0.433	ng	0.00

Target Compounds Qvalue

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

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