

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092421\
 Data File : BN016436.D
 Acq On : 24 Sep 2021 12:22
 Operator : CG/JU
 Sample : PB139115BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB139115BL

Quant Time: Sep 24 12:59:00 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	23732	0.40	ng	0.00
7) Naphthalene-d8	10.432	136	103173	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	50437	0.40	ng	0.00
19) Phenanthrene-d10	17.042	188	95945	0.40	ng	0.00
29) Chrysene-d12	21.259	240	73974	0.40	ng	# 0.01
35) Perylene-d12	23.518	264	60477	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.309	112	15444	0.29	ng	0.03
5) Phenol-d6	6.849	99	16269	0.26	ng	0.00
8) Nitrobenzene-d5	8.809	82	16864	0.23	ng	0.00
11) 2-Methylnaphthalene-d10	12.029	152	45520	0.30	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	2324	0.25	ng	0.00
15) 2-Fluorobiphenyl	12.911	172	60188	0.28	ng	-0.01
27) Fluoranthene-d10	19.087	212	72006	0.27	ng	0.00
31) Terphenyl-d14	19.698	244	58817	0.34	ng	0.00

Target Compounds Qvalue

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

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