

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100622\
 Data File : BN022016.D
 Acq On : 06 Oct 2022 11:20
 Operator : CG/JU
 Sample : PB147964BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147964BL

Quant Time: Oct 07 00:39:36 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	5921	0.400	ng	0.00
7) Naphthalene-d8	10.816	136	18118	0.400	ng	0.00
13) Acenaphthene-d10	14.653	164	9234	0.400	ng	0.00
19) Phenanthrene-d10	17.400	188	18312	0.400	ng	#-0.01
29) Chrysene-d12	21.600	240	13679	0.400	ng	# 0.00
35) Perylene-d12	24.081	264	11832	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.516	112	5661	0.413	ng	0.00
5) Phenol-d6	7.141	99	6936	0.390	ng	0.00
8) Nitrobenzene-d5	9.161	82	4877	0.335	ng	0.00
11) 2-Methylnaphthalene-d10	12.410	152	12309	0.370	ng	0.00
14) 2,4,6-Tribromophenol	16.146	330	895	0.254	ng	0.00
15) 2-Fluorobiphenyl	13.274	172	13871	0.344	ng	-0.01
27) Fluoranthene-d10	19.441	212	17433	0.358	ng	0.00
31) Terphenyl-d14	20.033	244	10323	0.375	ng	0.00

Target Compounds Qvalue

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

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