

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091522\
 Data File : BN021767.D
 Acq On : 15 Sep 2022 11:19
 Operator : CG/JU
 Sample : PB147570BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147570BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/16/2022
 Supervised By : Jagrut Upadhyay 09/16/2022

Quant Time: Sep 16 06:23:27 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 14 01:55:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.054	152	5312	0.400	ng	0.00
7) Naphthalene-d8	10.887	136	15882	0.400	ng	0.00
13) Acenaphthene-d10	14.707	164	8675	0.400	ng	0.00
19) Phenanthrene-d10	17.456	188	19702	0.400	ng	0.00
29) Chrysene-d12	21.647	240	14775	0.400	ng	# 0.00
35) Perylene-d12	24.160	264	12573	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.584	112	5858	0.422	ng	0.00
5) Phenol-d6	7.216	99	7062	0.401	ng	0.00
8) Nitrobenzene-d5	9.232	82	4886m	0.372	ng	0.00
11) 2-Methylnaphthalene-d10	12.474	152	14954	0.373	ng	0.00
14) 2,4,6-Tribromophenol	16.208	330	1147	0.289	ng	0.00
15) 2-Fluorobiphenyl	13.339	172	13315	0.376	ng	0.00
27) Fluoranthene-d10	19.490	212	21279	0.401	ng	0.00
31) Terphenyl-d14	20.080	244	13969	0.414	ng	0.00

Target Compounds Qvalue

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

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