

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092922\
 Data File : BN021953.D
 Acq On : 29 Sep 2022 12:29
 Operator : CG/JU
 Sample : PB147964BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147964BL

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 09/30/2022
 Supervised By :Jagrut Upadhyay 09/30/2022

Quant Time: Sep 30 02:44:39 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN092922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 30 02:27:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.025	152	4068	0.400	ng	0.00
7) Naphthalene-d8	10.866	136	11613	0.400	ng	# 0.01
13) Acenaphthene-d10	14.686	164	6022	0.400	ng	0.00
19) Phenanthrene-d10	17.437	188	11760	0.400	ng	# 0.00
29) Chrysene-d12	21.636	240	7834	0.400	ng	# 0.00
35) Perylene-d12	24.132	264	6428	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.563	112	3698	0.342	ng	0.00
5) Phenol-d6	7.195	99	4104	0.300	ng	0.00
8) Nitrobenzene-d5	9.211	82	3018m	0.311	ng	0.00
11) 2-Methylnaphthalene-d10	12.456	152	9094	0.363	ng	0.00
14) 2,4,6-Tribromophenol	16.196	330	585	0.227	ng	0.01
15) 2-Fluorobiphenyl	13.317	172	8716	0.327	ng	0.00
27) Fluoranthene-d10	19.475	212	10573	0.332	ng	0.00
31) Terphenyl-d14	20.060	244	6441	0.372	ng	0.00

Target Compounds Qvalue

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

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