

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091222\
 Data File : BN021716.D
 Acq On : 12 Sep 2022 11:23
 Operator : CG/JU
 Sample : PB147570BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147570BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/13/2022
 Supervised By : Jagrut Upadhyay 09/13/2022

Quant Time: Sep 13 02:33:49 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN082922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 13 02:32:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.062	152	6336	0.400	ng	0.00
7) Naphthalene-d8	10.888	136	19408	0.400	ng	0.00
13) Acenaphthene-d10	14.718	164	10489	0.400	ng	0.01
19) Phenanthrene-d10	17.470	188	20375	0.400	ng	0.01
29) Chrysene-d12	21.655	240	12947	0.400	ng	# 0.00
35) Perylene-d12	24.168	264	10859	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.592	112	6646	0.536	ng	0.00
5) Phenol-d6	7.217	99	7892	0.499	ng	0.00
8) Nitrobenzene-d5	9.233	82	5736m	0.438	ng	0.00
11) 2-Methylnaphthalene-d10	12.478	152	19274	0.441	ng	0.00
14) 2,4,6-Tribromophenol	16.219	330	1138	0.330	ng	0.01
15) 2-Fluorobiphenyl	13.339	172	16338	0.356	ng	0.00
27) Fluoranthene-d10	19.497	212	19954	0.381	ng	0.00
31) Terphenyl-d14	20.088	244	12438	0.427	ng	0.00

Target Compounds Qvalue

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

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