

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091422\
 Data File : BN021738.D
 Acq On : 13 Sep 2022 23:43
 Operator : CG/JU
 Sample : PB147574BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147574BL

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 09/15/2022
 Supervised By :Jagrut Upadhyay 09/15/2022

Quant Time: Sep 14 02:12:37 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	4602	0.400	ng	0.00
7) Naphthalene-d8	10.887	136	13314	0.400	ng	0.00
13) Acenaphthene-d10	14.707	164	7156	0.400	ng	0.00
19) Phenanthrene-d10	17.456	188	16810	0.400	ng	0.00
29) Chrysene-d12	21.647	240	11334	0.400	ng	# 0.00
35) Perylene-d12	24.160	264	9203	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.584	112	4108	0.341	ng	0.00
5) Phenol-d6	7.209	99	4958	0.325	ng	0.00
8) Nitrobenzene-d5	9.233	82	3427m	0.311	ng	0.00
11) 2-Methylnaphthalene-d10	12.475	152	11314	0.337	ng	0.00
14) 2,4,6-Tribromophenol	16.209	330	702	0.215	ng	0.00
15) 2-Fluorobiphenyl	13.339	172	10535	0.361	ng	0.00
27) Fluoranthene-d10	19.494	212	15882	0.351	ng	0.00
31) Terphenyl-d14	20.080	244	10379	0.401	ng	0.00

Target Compounds Qvalue

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

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