

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090122\
 Data File : BN021529.D
 Acq On : 01 Sep 2022 17:24
 Operator : CG/JU
 Sample : PB147332BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147332BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/02/2022
 Supervised By : Jagrut Upadhyay 09/02/2022

Quant Time: Sep 02 00:59:00 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN082922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 02 00:53:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.069	152	4333	0.400	ng	0.00
7) Naphthalene-d8	10.898	136	13312	0.400	ng	0.00
13) Acenaphthene-d10	14.729	164	6883	0.400	ng	0.01
19) Phenanthrene-d10	17.480	188	14077	0.400	ng	0.01
29) Chrysene-d12	21.664	240	8693	0.400	ng	# 0.00
35) Perylene-d12	24.191	264	6905	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.599	112	3533	0.417	ng	0.00
5) Phenol-d6	7.224	99	4228	0.391	ng	0.00
8) Nitrobenzene-d5	9.243	82	3012m	0.335	ng	0.00
11) 2-Methylnaphthalene-d10	12.490	152	12578	0.419	ng	0.00
14) 2,4,6-Tribromophenol	16.219	330	432	0.191	ng	0.00
15) 2-Fluorobiphenyl	13.349	172	10682	0.355	ng	0.00
27) Fluoranthene-d10	19.506	212	13364	0.369	ng	0.00
31) Terphenyl-d14	20.097	244	8583	0.439	ng	0.00

Target Compounds Qvalue

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

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