

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030722\
 Data File : BN018857.D
 Acq On : 07 Mar 2022 13:34
 Operator : CG/JU
 Sample : PB143119BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143119BL

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 03/08/2022
 Supervised By :Jagrut Upadhyay 03/08/2022

Quant Time: Mar 08 02:20:55 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN030322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 03 00:44:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.912	152	4386	0.400	ng	0.00
7) Naphthalene-d8	10.722	136	11970	0.400	ng	0.00
13) Acenaphthene-d10	14.538	164	7566	0.400	ng	0.00
19) Phenanthrene-d10	17.277	188	19503	0.400	ng	0.00
29) Chrysene-d12	21.458	240	20292	0.400	ng	0.00
35) Perylene-d12	23.849	264	17206	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.464	112	3398	0.354	ng	0.00
5) Phenol-d6	7.067	99	4104	0.357	ng	0.00
8) Nitrobenzene-d5	9.067	82	2649m	0.304	ng	0.00
11) 2-Methylnaphthalene-d10	12.303	152	5050	0.259	ng	0.00
14) 2,4,6-Tribromophenol	16.025	330	1054	0.302	ng	0.00
15) 2-Fluorobiphenyl	13.169	172	8684	0.268	ng	0.00
27) Fluoranthene-d10	19.307	212	15678	0.273	ng	0.00
31) Terphenyl-d14	19.898	244	13458	0.310	ng	0.00

Target Compounds Qvalue

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

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