

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040622\
 Data File : BN019367.D
 Acq On : 06 Apr 2022 12:47
 Operator : CG/JU
 Sample : PB143891BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143891BL

Quant Time: Apr 07 05:28:01 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN040422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 04 17:40:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.826	152	2977	0.400	ng	0.00
7) Naphthalene-d8	10.626	136	9350	0.400	ng	# 0.00
13) Acenaphthene-d10	14.463	164	6748	0.400	ng	0.00
19) Phenanthrene-d10	17.205	188	15561	0.400	ng	0.00
29) Chrysene-d12	21.396	240	14358	0.400	ng	0.00
35) Perylene-d12	23.741	264	13244	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.385	112	3131	0.393	ng	0.00
5) Phenol-d6	6.995	99	3517	0.367	ng	0.00
8) Nitrobenzene-d5	8.982	82	2304	0.271	ng	0.00
11) 2-Methylnaphthalene-d10	12.223	152	5227	0.318	ng	0.00
14) 2,4,6-Tribromophenol	15.954	330	1215	0.296	ng	0.00
15) 2-Fluorobiphenyl	13.095	172	7838	0.279	ng	0.00
27) Fluoranthene-d10	19.236	212	16478	0.329	ng	0.00
31) Terphenyl-d14	19.830	244	10644	0.341	ng	0.00

Target Compounds Qvalue

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

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