

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040822\
 Data File : BN019391.D
 Acq On : 08 Apr 2022 16:46
 Operator : CG/JU
 Sample : PB143983BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143983BL

Quant Time: Apr 11 04:10:57 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	7.818	152	5652	0.400	ng	0.00
7) Naphthalene-d8	10.626	136	16245	0.400	ng	# 0.00
13) Acenaphthene-d10	14.463	164	9627	0.400	ng	0.00
19) Phenanthrene-d10	17.205	188	22180	0.400	ng	# 0.00
29) Chrysene-d12	21.395	240	19547	0.400	ng	# 0.00
35) Perylene-d12	23.749	264	16969	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.370	112	5988	0.396	ng	-0.01
5) Phenol-d6	6.980	99	6707	0.368	ng	0.00
8) Nitrobenzene-d5	8.971	82	4289	0.291	ng	-0.01
11) 2-Methylnaphthalene-d10	12.216	152	8620	0.302	ng	0.00
14) 2,4,6-Tribromophenol	15.954	330	1627	0.278	ng	0.00
15) 2-Fluorobiphenyl	13.084	172	12015	0.300	ng	-0.01
27) Fluoranthene-d10	19.236	212	23563	0.330	ng	0.00
31) Terphenyl-d14	19.834	244	14323	0.337	ng	0.00

Target Compounds Qvalue

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

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