

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082622\
 Data File : BN021399.D
 Acq On : 27 Aug 2022 01:56
 Operator : CG/JU
 Sample : PB147294BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147294BL

Quant Time: Aug 27 04:06:59 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN082422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 23:37:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.083	152	4814	0.400	ng	0.00
7) Naphthalene-d8	10.909	136	14648	0.400	ng	0.00
13) Acenaphthene-d10	14.728	164	7724	0.400	ng	0.00
19) Phenanthrene-d10	17.480	188	16446	0.400	ng	0.00
29) Chrysene-d12	21.664	240	11492	0.400	ng	# 0.00
35) Perylene-d12	24.194	264	9131	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.599	112	3129	0.240	ng	0.00
5) Phenol-d6	7.231	99	4029	0.247	ng	0.00
8) Nitrobenzene-d5	9.254	82	2912	0.263	ng	0.00
11) 2-Methylnaphthalene-d10	12.497	152	9795	0.396	ng	0.00
14) 2,4,6-Tribromophenol	16.229	330	698	0.197	ng	0.01
15) 2-Fluorobiphenyl	13.360	172	11295	0.327	ng	0.00
27) Fluoranthene-d10	19.510	212	14815	0.301	ng	0.00
31) Terphenyl-d14	20.097	244	11004	0.453	ng	0.00

Target Compounds Qvalue

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

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