

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091022\
 Data File : BN021682.D
 Acq On : 10 Sep 2022 04:51
 Operator : CG/JU
 Sample : PB147497BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147497BL

Quant Time: Sep 10 05:28:31 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 09 23:19:14 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.062	152	4371	0.400	ng	0.00
7) Naphthalene-d8	10.898	136	13265	0.400	ng	0.00
13) Acenaphthene-d10	14.718	164	7253	0.400	ng	0.00
19) Phenanthrene-d10	17.470	188	14385	0.400	ng	0.00
29) Chrysene-d12	21.655	240	8909	0.400	ng	# 0.00
35) Perylene-d12	24.165	264	7585	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.592	112	5422	0.634	ng	0.00
5) Phenol-d6	7.217	99	6344	0.582	ng	0.00
8) Nitrobenzene-d5	9.233	82	4670	0.521	ng	0.00
11) 2-Methylnaphthalene-d10	12.482	152	14795	0.495	ng	0.00
14) 2,4,6-Tribromophenol	16.218	330	1122	0.471	ng	0.01
15) 2-Fluorobiphenyl	13.349	172	13106	0.413	ng	0.01
27) Fluoranthene-d10	19.493	212	17065	0.461	ng	0.00
31) Terphenyl-d14	20.088	244	10596	0.529	ng	0.00

Target Compounds Qvalue

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

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