

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081822\
 Data File : BF129887.D
 Acq On : 18 Aug 2022 21:19
 Operator : CG\JU
 Sample : N4200-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S2

Quant Time: Aug 19 01:04:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 18 12:03:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.793	152	112512	20.000	ng	0.00
21) Naphthalene-d8	8.075	136	450980	20.000	ng	0.00
39) Acenaphthene-d10	9.828	164	231888	20.000	ng	0.00
64) Phenanthrene-d10	11.322	188	361377	20.000	ng	0.00
76) Chrysene-d12	13.969	240	233628	20.000	ng	0.00
86) Perylene-d12	15.427	264	252356	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	878319	129.251	ng	0.01
7) Phenol-d6	6.451	99	996331	117.081	ng	0.00
23) Nitrobenzene-d5	7.363	82	757353	89.899	ng	0.00
42) 2,4,6-Tribromophenol	10.628	330	311730	133.942	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	1376760	90.174	ng	0.00
79) Terphenyl-d14	12.910	244	1257469	89.559	ng	0.00
Target Compounds						
77) Benzidine	12.663	184	111820	16.556	ng	Qvalue 100

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

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