

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040822\
 Data File : BF127723.D
 Acq On : 08 Apr 2022 19:54
 Operator : CG\JU
 Sample : N2207-06RX
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220350-AMENDED-096-21-02RX

Quant Time: Apr 11 02:23:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 06 13:55:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	146088	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	544917	20.000	ng	# 0.00
39) Acenaphthene-d10	10.010	164	318866	20.000	ng	0.00
64) Phenanthrene-d10	11.504	188	565853	20.000	ng	0.00
76) Chrysene-d12	14.151	240	352339	20.000	ng	0.00
86) Perylene-d12	15.680	264	304309	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	12609	1.402	ng	0.00
7) Phenol-d6	6.575	99	235324	20.008	ng	0.00
23) Nitrobenzene-d5	7.528	82	662136	54.171	ng	0.00
42) 2,4,6-Tribromophenol	10.798	330	6454	1.902	ng	0.00
45) 2-Fluorobiphenyl	9.328	172	1150630	51.048	ng	0.00
79) Terphenyl-d14	13.092	244	1152925	54.011	ng	0.00
Target Compounds						
77) Benzidine	12.839	184	85875	9.338	ng	Qvalue 99

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

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