

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092223\
 Data File : BF135407.D
 Acq On : 22 Sep 2023 19:09
 Operator : CG\JU
 Sample : 04434-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DIRTY-PILE-3

Quant Time: Sep 23 04:57:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 00:19:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.757	152	196193	20.000	ng	0.00
21) Naphthalene-d8	8.033	136	720826	20.000	ng	#-0.01
39) Acenaphthene-d10	9.792	164	347388	20.000	ng	-0.01
64) Phenanthrene-d10	11.286	188	558756	20.000	ng	# 0.00
76) Chrysene-d12	14.021	240	89051	20.000	ng	# 0.08
86) Perylene-d12	15.562	264	100810	20.000	ng	# 0.15
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	656840	54.452	ng	-0.01
7) Phenol-d6	6.398	99	803277	52.370	ng	0.00
23) Nitrobenzene-d5	7.322	82	475115	35.810	ng	-0.01
42) 2,4,6-Tribromophenol	10.586	330	137729	39.896	ng	-0.01
45) 2-Fluorobiphenyl	9.116	172	776278	33.714	ng	-0.01
79) Terphenyl-d14	12.910	244	275715	47.996	ng	0.02

Target Compounds	Qvalue

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

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