

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060922\
 Data File : BF128749.D
 Acq On : 09 Jun 2022 16:20
 Operator : CG\JU
 Sample : N3268-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-215(36-38)

Quant Time: Jun 10 03:28:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 08 14:28:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.810	152	77746	20.000	ng	0.00
21) Naphthalene-d8	8.098	136	299241	20.000	ng	# 0.00
39) Acenaphthene-d10	9.851	164	170537	20.000	ng	0.00
64) Phenanthrene-d10	11.339	188	304895	20.000	ng	0.00
76) Chrysene-d12	13.986	240	183675	20.000	ng	# 0.00
86) Perylene-d12	15.445	264	169914	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	682793	139.888	ng	0.00
7) Phenol-d6	6.475	99	878967	147.525	ng	0.00
23) Nitrobenzene-d5	7.381	82	617781	96.120	ng	0.00
42) 2,4,6-Tribromophenol	10.651	330	283668	140.780	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	1149219	94.397	ng	0.00
79) Terphenyl-d14	12.933	244	1122129	106.158	ng	0.00

Target Compounds	Qvalue

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

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