

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121621\
 Data File : BF126218.D
 Acq On : 16 Dec 2021 18:39
 Operator : JU/CG
 Sample : M4927-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PH1-BOT-14

Quant Time: Dec 17 01:47:15 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 12:22:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.816	152	183748	20.000	ng	0.00
21) Naphthalene-d8	8.098	136	756536	20.000	ng	# 0.00
39) Acenaphthene-d10	9.851	164	332882	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	460504	20.000	ng	# 0.00
76) Chrysene-d12	13.968	240	346285	20.000	ng	# 0.00
86) Perylene-d12	15.409	264	268842	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1354238	108.535	ng	0.00
7) Phenol-d6	6.439	99	1686136	106.427	ng	0.00
23) Nitrobenzene-d5	7.375	82	1065251	76.260	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	263843	98.626	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	1441688	75.959	ng	0.00
79) Terphenyl-d14	12.921	244	1127093	56.572	ng	0.00
Target Compounds						
10) Phenol	6.445	94	53256	2.995	ng	Qvalue 87

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

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