

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122821\
 Data File : BF126382.D
 Acq On : 28 Dec 2021 15:33
 Operator : JU/CG
 Sample : M5187-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2-119-1-5-10

Quant Time: Dec 28 16:09:45 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 13:00:11 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.810	152	111418	20.000	ng	0.00
21) Naphthalene-d8	8.092	136	456780	20.000	ng	#-0.01
39) Acenaphthene-d10	9.845	164	203092	20.000	ng	-0.01
64) Phenanthrene-d10	11.333	188	301073	20.000	ng	# 0.00
76) Chrysene-d12	13.968	240	211242	20.000	ng	#-0.01
86) Perylene-d12	15.421	264	232835	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	644635	85.204	ng	0.00
7) Phenol-d6	6.439	99	831375	86.541	ng	0.00
23) Nitrobenzene-d5	7.369	82	533149	63.214	ng	-0.01
42) 2,4,6-Tribromophenol	10.633	330	136186	83.440	ng	-0.01
45) 2-Fluorobiphenyl	9.169	172	748778	64.664	ng	-0.01
79) Terphenyl-d14	12.921	244	569567	46.864	ng	0.00

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

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