

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111321\
 Data File : BF126183.D
 Acq On : 13 Nov 2021 23:07
 Operator : JU/CG
 Sample : M4573-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MBR2

Quant Time: Nov 15 10:03:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 08:09:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	266516	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	1001620	20.000	ng	# 0.00
39) Acenaphthene-d10	9.869	164	574928	20.000	ng	0.00
64) Phenanthrene-d10	11.357	188	910345	20.000	ng	0.00
76) Chrysene-d12	13.992	240	506039	20.000	ng	#-0.01
86) Perylene-d12	15.456	264	628363	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.445	112	735377	47.384	ng	0.00
7) Phenol-d6	6.451	99	894909	40.905	ng	-0.02
23) Nitrobenzene-d5	7.392	82	1843233	86.787	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	716855	108.801	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	3254694	76.501	ng	0.00
79) Terphenyl-d14	12.945	244	2151043	68.545	ng	0.00
Target Compounds						
32) Benzoic acid	7.827	122	2024	9.331	ng	# 77

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

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