

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
 Data File : BF126141.D
 Acq On : 11 Nov 2021 23:35
 Operator : JU/CG
 Sample : M4639-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TT-3/4-5

Quant Time: Nov 12 09:44:52 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 03 17:21:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.839	152	145330	20.000	ng	-0.02
21) Naphthalene-d8	8.122	136	523680	20.000	ng	#-0.02
39) Acenaphthene-d10	9.880	164	272862	20.000	ng	-0.01
64) Phenanthrene-d10	11.363	188	436592	20.000	ng	-0.01
76) Chrysene-d12	14.004	240	414533	20.000	ng	# 0.00
86) Perylene-d12	15.468	264	266603	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	731906	86.485	ng	0.00
7) Phenol-d6	6.469	99	947221	79.400	ng	-0.02
23) Nitrobenzene-d5	7.404	82	691935	62.313	ng	-0.02
42) 2,4,6-Tribromophenol	10.669	330	244984	78.344	ng	-0.01
45) 2-Fluorobiphenyl	9.198	172	1240758	61.449	ng	-0.02
79) Terphenyl-d14	12.951	244	1245638	48.456	ng	-0.01

Target Compounds	Qvalue

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

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