

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121621\
 Data File : BF126217.D
 Acq On : 16 Dec 2021 18:06
 Operator : JU/CG
 Sample : M5022-09RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP10RE

Quant Time: Dec 17 01:47:03 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	192258	20.000	ng	0.00
21) Naphthalene-d8	8.098	136	812060	20.000	ng	# 0.00
39) Acenaphthene-d10	9.851	164	361252	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	498401	20.000	ng	# 0.00
76) Chrysene-d12	13.968	240	382453	20.000	ng	0.00
86) Perylene-d12	15.409	264	285991	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1327026	101.647	ng	0.00
7) Phenol-d6	6.439	99	1636498	98.722	ng	0.00
23) Nitrobenzene-d5	7.375	82	1039910	69.355	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	255559	88.027	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	1380968	67.046	ng	0.00
79) Terphenyl-d14	12.921	244	1084953	49.307	ng	0.00
Target Compounds						
10) Phenol	6.445	94	37443	2.012	ng	Qvalue 78

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

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