

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121621\
 Data File : BP008452.D
 Acq On : 16 Dec 2021 14:49
 Operator : CG/JU
 Sample : M5067-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BASE

Quant Time: Dec 16 15:51:05 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 14 17:59:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.734	152	51559	20.000	ng	-0.01
21) Naphthalene-d8	10.528	136	178359	20.000	ng	-0.01
39) Acenaphthene-d10	14.392	164	111581	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	239074	20.000	ng	0.00
76) Chrysene-d12	21.269	240	275424	20.000	ng	# 0.00
86) Perylene-d12	23.639	264	331413	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	398191	121.965	ng	0.00
7) Phenol-d6	6.934	99	436868	99.632	ng	0.00
23) Nitrobenzene-d5	8.899	82	374179	88.877	ng	-0.01
42) 2,4,6-Tribromophenol	15.898	330	245149	149.502	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	738371	84.120	ng	0.00
79) Terphenyl-d14	19.733	244	1331474	84.135	ng	0.00
Target Compounds						
84) Bis(2-ethylhexyl)phtha...	21.174	149	23314	2.308	ng	# 99

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

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