

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121621\
 Data File : BP008461.D
 Acq On : 16 Dec 2021 19:55
 Operator : CG/JU
 Sample : M5067-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BRICK

Quant Time: Dec 17 03:23:50 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	46047	20.000	ng	-0.01
21) Naphthalene-d8	10.534	136	163386	20.000	ng	0.00
39) Acenaphthene-d10	14.393	164	100442	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	207350	20.000	ng	0.00
76) Chrysene-d12	21.263	240	248944	20.000	ng	0.00
86) Perylene-d12	23.639	264	304243	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	347097	119.041	ng	-0.01
7) Phenol-d6	6.934	99	375318	95.841	ng	0.00
23) Nitrobenzene-d5	8.899	82	320263	83.042	ng	-0.01
42) 2,4,6-Tribromophenol	15.898	330	191464	129.712	ng	0.00
45) 2-Fluorobiphenyl	13.004	172	648942	82.130	ng	-0.01
79) Terphenyl-d14	19.728	244	1140934	79.764	ng	0.00

Target Compounds	Qvalue

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

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