

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120921\
 Data File : BP008323.D
 Acq On : 09 Dec 2021 21:41
 Operator : CG/JU
 Sample : M4967-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-7D

Quant Time: Dec 10 04:11:54 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 09 01:43:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	74776	20.000	ng	0.00
21) Naphthalene-d8	10.563	136	274861	20.000	ng	0.00
39) Acenaphthene-d10	14.422	164	175498	20.000	ng	0.00
64) Phenanthrene-d10	17.186	188	387861	20.000	ng	0.00
76) Chrysene-d12	21.292	240	436834	20.000	ng	0.00
86) Perylene-d12	23.674	264	476065	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	650050	139.970	ng	0.00
7) Phenol-d6	6.957	99	856841	136.672	ng	0.00
23) Nitrobenzene-d5	8.928	82	584795	96.749	ng	-0.01
42) 2,4,6-Tribromophenol	15.922	330	316997	141.317	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	1100483	91.083	ng	0.00
79) Terphenyl-d14	19.757	244	1838078	87.938	ng	-0.01
Target Compounds						
10) Phenol	6.981	94	13658	2.120	ng	Qvalue # 50
84) Bis(2-ethylhexyl)phtha...	21.204	149	56873	3.011	ng	99

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

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