

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010322\
 Data File : BP008634.D
 Acq On : 03 Jan 2022 13:07
 Operator : CG/JU
 Sample : PB141757TB
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB141757TB

Quant Time: Jan 03 14:11:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 30 08:44:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	630881	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	2444452	20.000	ng	0.00
39) Acenaphthene-d10	14.522	164	1644486	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	3568769	20.000	ng	0.00
76) Chrysene-d12	21.357	240	3233858	20.000	ng	0.00
86) Perylene-d12	23.780	264	3046027	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	4773108	138.096	ng	0.00
7) Phenol-d6	7.063	99	5812464	120.817	ng	0.00
23) Nitrobenzene-d5	9.046	82	3451678	84.001	ng	-0.01
42) 2,4,6-Tribromophenol	16.016	330	3262753	146.646	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	9652485	85.040	ng	0.00
79) Terphenyl-d14	19.827	244	13277550	83.832	ng	0.00

Target Compounds	Qvalue

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

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