

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120121\
 Data File : BP008168.D
 Acq On : 01 Dec 2021 11:19
 Operator : CG/JU
 Sample : PB141067BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB141067BL

Quant Time: Dec 01 12:07:09 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 01 12:06:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	73910	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	260973	20.000	ng	0.00
39) Acenaphthene-d10	14.445	164	175199	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	401478	20.000	ng	0.00
76) Chrysene-d12	21.304	240	457596	20.000	ng	0.00
86) Perylene-d12	23.698	264	488749	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	633438	137.991	ng	0.00
7) Phenol-d6	6.987	99	796650	128.560	ng	0.00
23) Nitrobenzene-d5	8.963	82	540152	94.119	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	346833	154.882	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	1100167	91.212	ng	0.00
79) Terphenyl-d14	19.780	244	2086324	95.286	ng	0.00

Target Compounds	Qvalue

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

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