

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110821\
 Data File : BP007871.D
 Acq On : 08 Nov 2021 20:57
 Operator : CG/JU
 Sample : M4576-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-2-(9-9.5)

Quant Time: Nov 09 08:45:38 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 08 16:36:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	104230	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	387983	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	252383	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	564722	20.000	ng	-0.01
76) Chrysene-d12	21.345	240	672613	20.000	ng	0.00
86) Perylene-d12	23.750	264	733331	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	850791	138.433	ng	0.00
7) Phenol-d6	7.040	99	1035213	116.375	ng	0.00
23) Nitrobenzene-d5	9.016	82	595531	82.453	ng	-0.02
42) 2,4,6-Tribromophenol	15.992	330	424919	117.972	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	1425990	82.281	ng	-0.01
79) Terphenyl-d14	19.815	244	2577493	78.032	ng	0.00

Target Compounds	Qvalue

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

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