

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
 Data File : BP007923.D
 Acq On : 10 Nov 2021 17:08
 Operator : CG/JU
 Sample : M4567-01RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BASEMENTRE

Quant Time: Nov 11 10:14:35 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 10:45:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	127667	20.000	ng	-0.01
21) Naphthalene-d8	10.652	136	425863	20.000	ng	-0.01
39) Acenaphthene-d10	14.493	164	269790	20.000	ng	-0.01
64) Phenanthrene-d10	17.245	188	637107	20.000	ng	-0.01
76) Chrysene-d12	21.339	240	739334	20.000	ng	-0.01
86) Perylene-d12	23.751	264	856841	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	51473	6.838	ng	0.00
7) Phenol-d6	7.034	99	94454	8.669	ng	0.00
23) Nitrobenzene-d5	9.011	82	267144	33.697	ng	-0.02
42) 2,4,6-Tribromophenol	15.987	330	32210	8.366	ng	-0.01
45) 2-Fluorobiphenyl	13.116	172	466880	25.201	ng	-0.01
79) Terphenyl-d14	19.810	244	518804	14.289	ng	-0.01

Target Compounds	Qvalue

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

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