

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
 Data File : BP007917.D
 Acq On : 10 Nov 2021 13:36
 Operator : CG/JU
 Sample : M4567-01RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BASEMENTRE

Quant Time: Nov 10 14:56:59 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	7.857	152	137469	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	477537	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	281160	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	529489	20.000	ng	0.00
76) Chrysene-d12	21.345	240	529830	20.000	ng	# 0.00
86) Perylene-d12	23.757	264	648913	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	60944	7.519	ng	0.00
7) Phenol-d6	7.034	99	111050	9.465	ng	0.00
23) Nitrobenzene-d5	9.016	82	292250	32.875	ng	-0.01
42) 2,4,6-Tribromophenol	15.986	330	26628	6.636	ng	-0.01
45) 2-Fluorobiphenyl	13.122	172	504385	26.125	ng	0.00
79) Terphenyl-d14	19.816	244	383357	14.733	ng	0.00

Target Compounds	Qvalue

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

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