

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110821\
 Data File : BP007875.D
 Acq On : 08 Nov 2021 23:12
 Operator : CG/JU
 Sample : M4523-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FIELD-BLANK

Quant Time: Nov 09 08:47:05 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	100112	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	368485	20.000	ng	-0.01
39) Acenaphthene-d10	14.498	164	235813	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	528222	20.000	ng	#-0.01
76) Chrysene-d12	21.339	240	650403	20.000	ng	-0.01
86) Perylene-d12	23.750	264	723639	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	469256	79.494	ng	0.00
7) Phenol-d6	7.034	99	350205	40.988	ng	-0.01
23) Nitrobenzene-d5	9.016	82	638430	93.070	ng	-0.02
42) 2,4,6-Tribromophenol	15.992	330	464873	138.133	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	1495970	92.384	ng	-0.01
79) Terphenyl-d14	19.815	244	2856144	89.420	ng	0.00

Target Compounds	Qvalue

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

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