

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082421\
 Data File : BP006870.D
 Acq On : 25 Aug 2021 04:29
 Operator : CG/JU
 Sample : M3468-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SAMPLE-6-(203.30)

Quant Time: Aug 25 05:01:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 24 13:40:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	153307	20.000	ng	0.00
21) Naphthalene-d8	10.463	136	632159	20.000	ng	#-0.01
39) Acenaphthene-d10	14.328	164	357504	20.000	ng	-0.01
64) Phenanthrene-d10	17.086	188	716449	20.000	ng	#-0.01
76) Chrysene-d12	21.192	240	696133	20.000	ng	#-0.02
86) Perylene-d12	23.492	264	746412	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	1421498	142.416	ng	0.00
7) Phenol-d6	6.875	99	1586525	111.896	ng	0.00
23) Nitrobenzene-d5	8.846	82	1320666	96.429	ng	0.00
42) 2,4,6-Tribromophenol	15.828	330	629776	168.556	ng	-0.02
45) 2-Fluorobiphenyl	12.951	172	2344066	98.097	ng	0.00
79) Terphenyl-d14	19.669	244	3726623	107.269	ng	-0.02

Target Compounds	Qvalue

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

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