

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082421\
 Data File : BP006868.D
 Acq On : 25 Aug 2021 03:21
 Operator : CG/JU
 Sample : M3468-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SAMPLE-4-(177.00)

Quant Time: Aug 25 04:54:57 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	148990	20.000	ng	0.00
21) Naphthalene-d8	10.463	136	609087	20.000	ng	#-0.01
39) Acenaphthene-d10	14.328	164	340624	20.000	ng	-0.01
64) Phenanthrene-d10	17.086	188	684461	20.000	ng	#-0.01
76) Chrysene-d12	21.192	240	668587	20.000	ng	#-0.02
86) Perylene-d12	23.492	264	724007	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	1299506	133.966	ng	0.00
7) Phenol-d6	6.875	99	1369541	99.391	ng	0.00
23) Nitrobenzene-d5	8.846	82	1286754	97.512	ng	0.00
42) 2,4,6-Tribromophenol	15.828	330	612711	172.115	ng	-0.02
45) 2-Fluorobiphenyl	12.951	172	2279673	100.130	ng	0.00
79) Terphenyl-d14	19.669	244	3602371	107.965	ng	-0.02

Target Compounds	Qvalue

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

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