

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082421\
 Data File : BP006871.D
 Acq On : 25 Aug 2021 05:03
 Operator : CG/JU
 Sample : M3468-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SAMPLE-9-(322.00)

Quant Time: Aug 25 05:46:52 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	148086	20.000	ng	0.00
21) Naphthalene-d8	10.463	136	600766	20.000	ng	#-0.01
39) Acenaphthene-d10	14.328	164	328438	20.000	ng	-0.01
64) Phenanthrene-d10	17.086	188	665916	20.000	ng	#-0.01
76) Chrysene-d12	21.192	240	648170	20.000	ng	#-0.02
86) Perylene-d12	23.492	264	691199	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	1329271	137.871	ng	0.00
7) Phenol-d6	6.875	99	1462691	106.799	ng	0.00
23) Nitrobenzene-d5	8.846	82	1241832	95.411	ng	0.00
42) 2,4,6-Tribromophenol	15.828	330	556702	162.184	ng	-0.02
45) 2-Fluorobiphenyl	12.945	172	2168446	98.779	ng	-0.01
79) Terphenyl-d14	19.669	244	3355764	103.742	ng	-0.02

Target Compounds	Qvalue

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

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