

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
 Data File : BP006186.D
 Acq On : 12 Jul 2021 17:56
 Operator : CG/JU
 Sample : M3002-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-F

Quant Time: Jul 12 18:49:31 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 12 14:59:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	192105	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	769431	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	441762	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	817116	20.000	ng	# 0.00
76) Chrysene-d12	21.321	240	737013	20.000	ng	# 0.00
86) Perylene-d12	23.704	264	822739	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	1185484	106.181	ng	0.00
7) Phenol-d6	7.052	99	1468040	99.407	ng	0.00
23) Nitrobenzene-d5	9.028	82	810835	66.564	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	510274	97.020	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	1504174	50.337	ng	0.00
79) Terphenyl-d14	19.798	244	1771110	47.380	ng	0.00
Target Compounds						
32) Benzoic acid	10.046	122	6692	7.649	ng	Qvalue 90

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

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