

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090822\
 Data File : BP011654.D
 Acq On : 08 Sep 2022 11:54
 Operator : CG/JU
 Sample : PB147456TB
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147456TB

Quant Time: Sep 09 04:27:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP090222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 03 04:07:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	587413	20.000	ng	0.00
21) Naphthalene-d8	10.787	136	2316178	20.000	ng	# 0.00
39) Acenaphthene-d10	14.616	164	1270145	20.000	ng	0.00
64) Phenanthrene-d10	17.369	188	2459436	20.000	ng	0.00
76) Chrysene-d12	21.451	240	1901266	20.000	ng	0.00
86) Perylene-d12	23.927	264	1666371	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	5971808	179.801	ng	0.00
7) Phenol-d6	7.134	99	7292891	165.942	ng	0.00
23) Nitrobenzene-d5	9.140	82	4238734	112.336	ng	0.00
42) 2,4,6-Tribromophenol	16.110	330	2277784	158.723	ng	0.00
45) 2-Fluorobiphenyl	13.246	172	8979696	101.098	ng	0.00
79) Terphenyl-d14	19.922	244	11317822	124.681	ng	0.00

Target Compounds Qvalue

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

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