

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092822\
 Data File : BP011831.D
 Acq On : 28 Sep 2022 15:44
 Operator : CG/JU
 Sample : PB147907TB
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147907TB

Quant Time: Sep 29 02:26:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:03:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	384543	20.000	ng	0.00
21) Naphthalene-d8	10.716	136	1649684	20.000	ng	# 0.00
39) Acenaphthene-d10	14.551	164	981091	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	1986056	20.000	ng	0.00
76) Chrysene-d12	21.386	240	1589892	20.000	ng	-0.01
86) Perylene-d12	23.827	264	1313240	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	3114537	142.111	ng	0.00
7) Phenol-d6	7.069	99	3991354	123.169	ng	0.00
23) Nitrobenzene-d5	9.075	82	2179119	69.138	ng	0.00
42) 2,4,6-Tribromophenol	16.045	330	1047137	158.509	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	4853889	78.238	ng	0.00
79) Terphenyl-d14	19.857	244	6493017	93.903	ng	-0.01

Target Compounds	Qvalue

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

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