

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092221\
 Data File : BP007091.D
 Acq On : 22 Sep 2021 10:40
 Operator : CG/JU
 Sample : PB139216BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB139216BL

Quant Time: Sep 22 11:34:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 16:50:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	260121	20.000	ng	0.00
21) Naphthalene-d8	10.699	136	1026458	20.000	ng	# 0.00
39) Acenaphthene-d10	14.528	164	666405	20.000	ng	0.00
64) Phenanthrene-d10	17.281	188	1478153	20.000	ng	# 0.00
76) Chrysene-d12	21.363	240	1462470	20.000	ng	# 0.00
86) Perylene-d12	23.774	264	1486360	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.470	112	1942904	125.031	ng	0.00
7) Phenol-d6	7.064	99	2363541	111.285	ng	0.00
23) Nitrobenzene-d5	9.058	82	1335714	75.782	ng	-0.01
42) 2,4,6-Tribromophenol	16.016	330	987136	114.257	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	3445453	74.070	ng	0.00
79) Terphenyl-d14	19.834	244	5828862	76.375	ng	0.00

Target Compounds	Qvalue

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

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