

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092321\
 Data File : BP007137.D
 Acq On : 23 Sep 2021 22:11
 Operator : CG/JU
 Sample : M3832-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-B

Quant Time: Sep 24 01:02:38 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 23 11:17:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	283964	20.000	ng	0.00
21) Naphthalene-d8	10.687	136	1112977	20.000	ng	# 0.00
39) Acenaphthene-d10	14.516	164	696927	20.000	ng	-0.01
64) Phenanthrene-d10	17.269	188	1414059	20.000	ng	0.00
76) Chrysene-d12	21.351	240	1392912	20.000	ng	# 0.00
86) Perylene-d12	23.757	264	1502027	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.470	112	2249076	132.581	ng	0.00
7) Phenol-d6	7.064	99	2832574	122.171	ng	0.00
23) Nitrobenzene-d5	9.052	82	1649733	86.322	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	1337940	148.079	ng	0.00
45) 2-Fluorobiphenyl	13.146	172	4074038	83.747	ng	0.00
79) Terphenyl-d14	19.828	244	5827775	80.174	ng	0.00

Target Compounds	Qvalue

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

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