

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091721\
 Data File : BP007037.D
 Acq On : 17 Sep 2021 21:19
 Operator : CG/JU
 Sample : M3791-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SW-CB-03-(1.0)

Manual Integrations
 APPROVED

mohammad
 9/20/2021 4:21:42 PM

Quant Time: Sep 18 00:58:01 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 17 10:50:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	312995	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1218759	20.000	ng	# 0.00
39) Acenaphthene-d10	14.534	164	739386	20.000	ng	0.00
64) Phenanthrene-d10	17.280	188	1531436	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	1545645	20.000	ng	#-0.01
86) Perylene-d12	23.768	264	1632724	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	1152473	60.126	ng	0.00
7) Phenol-d6	7.069	99	850564	32.511	ng	0.00
23) Nitrobenzene-d5	9.069	82	1910814	88.057	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	1278015	143.727	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	4696894	87.696	ng	0.00
79) Terphenyl-d14	19.833	244	7147147	88.718	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	9.916	122	3429m	6.329	ng	

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

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