

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092021\
 Data File : BP007069.D
 Acq On : 21 Sep 2021 01:56
 Operator : CG/JU
 Sample : M3803-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 26160

Manual Integrations
 APPROVED

mohammad
 9/21/2021 11:29:47 AM

Quant Time: Sep 21 02:41:08 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	330217	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1307938	20.000	ng	# 0.00
39) Acenaphthene-d10	14.534	164	825499	20.000	ng	0.00
64) Phenanthrene-d10	17.281	188	1656445	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	1632108	20.000	ng	#-0.01
86) Perylene-d12	23.774	264	1744011	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	2356434	116.526	ng	0.00
7) Phenol-d6	7.075	99	3021521	109.469	ng	0.00
23) Nitrobenzene-d5	9.063	82	1710244	73.440	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	1248913	125.802	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	3888803	65.034	ng	-0.01
79) Terphenyl-d14	19.833	244	5752849	67.627	ng	0.00
Target Compounds						
32) Benzoic acid	9.934	122	4733m	6.395	ng	Qvalue
84) Bis(2-ethylhexyl)phtha...	21.269	149	131587	2.261	ng	95

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

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