Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092021\
Data File : BP007064.D
Acq On : 20 Sep 2021 23:07
Operator : CG/JU
Sample : M3817-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SW-CB-04-(1.5)

Quant Time: Sep 21 01:04:24 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	340229	20.000	ng	0.00
21) Naphthalene-d8	10.705	136	1316713	20.000	ng	# 0.00
39) Acenaphthene-d10	14.534	164	819040	20.000	ng	0.00
64) Phenanthrene-d10	17.281	188	1709042	20.000	ng	# 0.00
76) Chrysene-d12	21.357	240	1713024	20.000	ng	#-0.01
86) Perylene-d12	23.774	264	1780048	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	929512	44.612	ng	0.00
7) Phenol-d6	7.069	99	674147	23.705	ng	0.00
23) Nitrobenzene-d5	9.063	82	1742982	74.347	ng	0.00
42) 2,4,6-Tribromophenol	16.022	330	1239123	125.801	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	4222694	71.174	ng	0.00
79) Terphenyl-d14	19.833	244	6300335	70.565	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.093	94	62351	2.175	ng	97
20) 3+4-Methylphenols	8.669	107	831641	33.365	ng	86
31) Naphthalene	10.757	128	293704	4.026	ng	99
32) Benzoic acid	9.975	122	114633	13.300	ng	94

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

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