Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121721\
Data File : BP008477.D
Acq On : 17 Dec 2021 18:27
Operator : CG/JU
Sample : M5084-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C4

Quant Time: Dec 18 07:16:04 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 18 06:49:58 2021
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	45826	20.000	ng	0.00
21) Naphthalene-d8	10.522	136	171420	20.000	ng	0.00
39) Acenaphthene-d10	14.381	164	117607	20.000	ng	0.00
64) Phenanthrene-d10	17.145	188	251141	20.000	ng	0.00
76) Chrysene-d12	21.257	240	291575	20.000	ng	0.00
86) Perylene-d12	23.627	264	331579	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.328	112	185022	63.762	ng	0.00
7) Phenol-d6	6.934	99	135402	34.743	ng	0.00
23) Nitrobenzene-d5	8.893	82	356821	88.185	ng	0.00
42) 2,4,6-Tribromophenol	15.887	330	263585	152.509	ng	0.00
45) 2-Fluorobiphenyl	12.999	172	772923	83.544	ng	0.00
79) Terphenyl-d14	19.722	244	1301093	77.661	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.964	94	13989	3.535	ng	77
32) Benzoic acid	9.916	122	18502	10.487	ng	# 33
52) Acenaphthene	14.446	154	19736	3.114	ng	97
58) Fluorene	15.440	166	32557	3.896	ng	96

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

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