Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
Data File : BP019835.D
Acq On : 22 Feb 2024 21:12
Operator : MA/JU
Sample : P1521-04
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CG-MW-17-CH3

Quant Time: Feb 23 00:59:55 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 19 23:35:42 2024
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	78268	20.000	ng	0.00
21) Naphthalene-d8	10.693	136	293559	20.000	ng	0.00
39) Acenaphthene-d10	14.534	164	209134	20.000	ng	0.00
64) Phenanthrene-d10	17.286	188	479189	20.000	ng	0.00
76) Chrysene-d12	21.386	240	463626	20.000	ng	0.00
86) Perylene-d12	23.810	264	519174	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	307902	63.412	ng	0.00
7) Phenol-d6	7.069	99	270254	41.279	ng	0.00
23) Nitrobenzene-d5	9.075	82	540227	79.741	ng	0.00
42) 2,4,6-Tribromophenol	16.028	330	454480	148.839	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	1299189	76.797	ng	0.00
79) Terphenyl-d14	19.857	244	2063138	73.182	ng	0.00
Target Compounds						Qvalue
6) Aniline	7.234	93	18883	2.968	ng	94
10) Phenol	7.099	94	51560	7.902	ng	95
20) 3+4-Methylphenols	8.728	107	11386978	1899.242	ng	89
31) Naphthalene	10.746	128	142249	8.835	ng	97
32) Benzoic acid	10.022	122	55933	17.262	ng	97

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

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