Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100422\
Data File : BG055073.D
Acq On : 4 Oct 2022 21:42
Operator : CG/JU
Sample : N4936-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SASP2-T-2-(16-22)

Quant Time: Oct 05 00:34:35 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 04 07:12:22 2022
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.344	152	32278	20.000	ng	0.00
21) Naphthalene-d8	11.176	136	139566	20.000	ng	0.00
39) Acenaphthene-d10	14.948	164	96923	20.000	ng	0.00
64) Phenanthrene-d10	17.680	188	212316	20.000	ng	0.00
76) Chrysene-d12	21.975	240	205371	20.000	ng	0.00
86) Perylene-d12	25.436	264	221531	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.864	112	178989	123.167	ng	0.00
7) Phenol-d6	7.468	99	247013	113.445	ng	0.00
23) Nitrobenzene-d5	9.513	82	155785	66.698	ng	0.00
42) 2,4,6-Tribromophenol	16.423	330	120463	120.106	ng	0.00
45) 2-Fluorobiphenyl	13.579	172	391428	63.047	ng	0.00
79) Terphenyl-d14	20.259	244	616674	60.691	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.721	178	37109	3.327	ng	97
75) Fluoranthene	19.713	202	38886	2.795	ng	99
78) Pyrene	20.071	202	34231	2.507	ng	98

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

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