Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062219\
Data File : BM021093.D
Acq On : 22 Jun 2019 10:10
Operator : HP/JU
Sample : K3309-03DL 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
KY032MW006-190610DL

Manual Integrations APPROVED
mohammad
6/26/2019 1:57:33 PM

Quant Time: Jun 27 08:46:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM062019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 20 15:39:58 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	317400	20.00	ng		0.00
21) Naphthalene-d8	10.57	136	1338390	20.00	ng		0.00
39) Acenaphthene-d10	14.42	164	866383	20.00	ng		0.00
64) Phenanthrene-d10	17.16	188	1990868	20.00	ng		0.00
76) Chrysene-d12	21.34	240	1628274	20.00	ng		0.00
87) Perylene-d12	23.62	264	1654213	20.00	ng		-0.01
System Monitoring Compounds							
5) 2-Fluorophenol	5.37	112	220466	12.58	ng		0.00
7) Phenol-d6	6.95	99	213674	8.37	ng		0.00
23) Nitrobenzene-d5	8.95	82	401593	15.64	ng		0.00
42) 2,4,6-Tribromophenol	15.91	330	374969	23.06	ng		0.00
45) 2-Fluorobiphenyl	13.03	172	1162428	16.49	ng		0.00
79) Terphenyl-d14	19.79	244	1390527	15.95	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.32	88	5839773	884.271	ng		97
4) n-Nitrosodimethylamine	3.62	42	372683	50.966	ng	#	1
10) Phenol	6.98	94	936246	35.902	ng		99
11) bis(2-Chloroethyl)ether	7.21	93	1490021	72.168	ng		98
13) 1,4-Dichlorobenzene	7.81	146	274755	10.340	ng		99
14) 1,2-Dichlorobenzene	8.13	146	1305808	50.143	ng		99
17) 2-Methylphenol	8.23	107	1291920	69.355	ng		100
20) 3+4-Methylphenols	8.57	107	8627969	336.496	ng		99
22) Acetophenone	8.61	105	115199	3.378	ng	#	84
27) 2,4-Dimethylphenol	9.79	122	3708702m	204.135	ng		
31) Naphthalene	10.62	128	504294	7.120	ng		99

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

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