Quantitation Report (LSC Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100118\
Data File : BF109482.D
Acq On : 1 Oct 2018 19:32
Operator : JU/SJ
Sample : J5155-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
QNWP8-MW30D-GW

Manual Integrations APPROVED
Sohil
10/2/2018 1:01:35 PM

Quant Time: Oct 02 04:16:45 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Sep 22 13:32:18 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	101254	20.00	ng		0.00
21) Naphthalene-d8	7.99	136	368236	20.00	ng		0.00
38) Acenaphthene-d10	9.74	164	159201	20.00	ng		0.00
63) Phenanthrene-d10	11.22	188	257284	20.00	ng		0.00
75) Chrysene-d12	13.85	240	234909	20.00	ng		0.00
86) Perylene-d12	15.24	264	181062	20.00	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.33	112	264876	43.09	ng		0.01
7) Phenol-d6	6.35	99	203416	25.93	ng		0.00
23) Nitrobenzene-d5	7.27	82	448565	71.91	ng		-0.01
41) 2,4,6-Tribromophenol	10.53	330	155844	99.30	ng		0.00
44) 2-Fluorobiphenyl	9.06	172	785567	74.42	ng		-0.01
78) Terphenyl-d14	12.80	244	661682	69.13	ng		0.00
Target Compounds							Qvalue
10) Phenol	6.37	94	24070	2.709	ng		77
20) 3+4-Methylphenols	7.13	107	13606	2.037	ng	#	78
22) Acetophenone	7.11	105	31533	3.704	ng	#	89
27) 2,4-Dimethylphenol	7.65	122	29808m	6.090	ng		
37) 2-Methylnaphthalene	8.70	142	75834m	6.836	ng		
45) 1,1'-Biphenyl	9.16	154	155814	12.012	ng		99
48) Acenaphthylene	9.60	152	53273	3.408	ng		97
49) Dimethylphthalate	9.46	163	240532	20.773	ng		99
51) Acenaphthene	9.77	154	697576	73.807	ng		100
54) Dibenzofuran	9.95	168	533579	37.961	ng		99
57) Fluorene	10.28	166	247034	24.632	ng	#	96
70) Phenanthrene	11.24	178	88328	6.876	ng		99
72) Carbazole	11.45	167	218618	16.864	ng		99

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

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