Quantitation Report (LSC Reviewed)

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF101217\
Data File : BF099416.D
Acq On : 13 Oct 2017 1:10
Operator : SJ/JU
Sample : I5729-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 7W-(1-2)

Manual Integrations APPROVED
Sohil
10/13/2017 3:06:38 PM

Quant Time: Oct 13 05:01:13 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 11 16:25:21 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	111539	20.00	ng		0.00
21) Naphthalene-d8	8.13	136	433807	20.00	ng		0.00
38) Acenaphthene-d10	9.89	164	169035	20.00	ng		0.00
63) Phenanthrene-d10	11.37	188	244915	20.00	ng		0.00
75) Chrysene-d12	14.02	240	209077	20.00	ng		0.00
86) Perylene-d12	15.49	264	174200	20.00	ng		0.01
System Monitoring Compounds							
5) 2-Fluorophenol	5.48	112	677214	96.65	ng		0.01
7) Phenol-d6	6.49	99	827592	95.75	ng		0.00
23) Nitrobenzene-d5	7.42	82	500999	74.06	ng		0.00
41) 2,4,6-Tribromophenol	10.68	330	116059	83.91	ng		0.00
44) 2-Fluorobiphenyl	9.21	172	787225	77.75	ng		0.00
78) Terphenyl-d14	12.96	244	532885	57.96	ng		0.00
Target Compounds							Qvalue
10) Phenol	6.50	94	27635	2.859	ng		95
31) Naphthalene	8.16	128	49496	2.429	ng		100
37) 2-Methylnaphthalene	8.85	142	35401	2.673	ng		98
49) Dimethylphthalate	9.60	163	182972	14.342	ng		99
51) Acenaphthene	9.92	154	34476	3.259	ng		99
54) Dibenzofuran	10.09	168	56225	3.992	ng		97
57) Fluorene	10.43	166	38973	3.443	ng		97
70) Phenanthrene	11.40	178	729144	55.619	ng		99
71) Anthracene	11.45	178	121486	9.057	ng		99
72) Carbazole	11.61	167	77870	5.928	ng		100
73) Di-n-butylphthalate	11.93	149	41543	2.627	ng		97
74) Fluoranthene	12.60	202	985225	74.217	ng		98
77) Pyrene	12.83	202	871864	54.687	ng		99
80) Benzo(a)anthracene	14.01	228	421896	33.325	ng		99
82) Chrysene	14.04	228	365206	30.300	ng		97
85) Indeno(1,2,3-cd)pyrene	16.97	276	137661	14.278	ng		99
87) Benzo(b)fluoranthene	15.06	252	414694m	38.718	ng		
88) Benzo(k)fluoranthene	15.09	252	143280m	13.544	ng		
89) Benzo(a)pyrene	15.43	252	252716	25.705	ng	#	95
90) Dibenzo(a,h)anthracene	16.97	278	34532	4.389	ng	#	77
91) Benzo(g,h,i)perylene	17.43	276	133010	17.102	ng		95

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

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