Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
Data File : BF106879.D
Acq On : 27 Jun 2018 18:28
Operator : JU/SJ
Sample : J3711-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : QNWP8-265-GW

Quant Time: Jun 28 11:35:50 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 25 18:50:17 2018
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	43112	20.00	ng		0.00
21) Naphthalene-d8	8.24	136	59516	20.00	ng		-0.01
38) Acenaphthene-d10	10.00	164	80720	20.00	ng		0.00
63) Phenanthrene-d10	11.50	188	110793	20.00	ng		0.00
75) Chrysene-d12	14.15	240	120597	20.00	ng		-0.02
86) Perylene-d12	15.69	264	137089	20.00	ng		-0.03
System Monitoring Compounds							
5) 2-Fluorophenol	5.62	112	112373	44.14	ng		0.02
7) Phenol-d6	6.62	99	110922	34.89	ng		0.00
23) Nitrobenzene-d5	7.53	82	236159	220.52	ng		0.00
41) 2,4,6-Tribromophenol	10.80	330	112872	120.65	ng		0.00
44) 2-Fluorobiphenyl	9.32	172	426124	90.27	ng		0.00
78) Terphenyl-d14	13.08	244	321214	64.52	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	2.80	88	10818	8.265	ng		98
3) Pyridine	3.58	79	57420	16.175	ng		98
6) Aniline	6.64	93	188828	43.782	ng	#	74
10) Phenol	6.63	94	85095	23.452	ng	#	25
17) 2-Methylphenol	7.22	107	90242	37.762	ng		98
20) 3+4-Methylphenols	7.38	107	413882	149.632	ng		93
22) Acetophenone	7.38	105	66368	49.844	ng	#	3
27) 2,4-Dimethylphenol	7.95	122	639048	801.018	ng		99
31) Naphthalene	8.27	128	5587683	2008.117	ng		98
37) 2-Methylnaphthalene	8.97	142	724107	392.609	ng		97
45) 1,1'-Biphenyl	9.42	154	311729	48.458	ng		98
48) Acenaphthylene	9.87	152	63383	7.928	ng	#	90
49) Dimethylphthalate	9.72	163	29772	4.918	ng	#	95
51) Acenaphthene	10.04	154	468614	93.085	ng		98
54) Dibenzofuran	10.21	168	578657	83.943	ng		98
57) Fluorene	10.55	166	352340	64.411	ng		99
70) Phenanthrene	11.52	178	471297	88.196	ng		99
71) Anthracene	11.57	178	76014	13.936	ng		99
72) Carbazole	11.74	167	364778	71.432	ng		98
74) Fluoranthene	12.73	202	49164	8.347	ng		95
77) Pyrene	12.95	202	34174	4.297	ng		99

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

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