

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100218\
 Data File : BF109496.D
 Acq On : 2 Oct 2018 13:01
 Operator : JU/SJ
 Sample : J5155-04DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW9-GWDL

Manual Integrations
 APPROVED

Sohil
 10/3/2018 9:30:59 AM

Quant Time: Oct 02 13:59:39 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	98882	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	358785	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	193401	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	376864	20.00	ng	0.00
75) Chrysene-d12	13.84	240	282961	20.00	ng	0.00
86) Perylene-d12	15.24	264	268665	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	6025	1.00	ng	0.01
7) Phenol-d6	6.36	99	16976	2.22	ng	0.00
23) Nitrobenzene-d5	7.26	82	27303	4.49	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	8619	4.52	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	51761	4.04	ng	-0.02
78) Terphenyl-d14	12.79	244	32618	2.83	ng	-0.02

Target Compounds

						Qvalue
6) Aniline	6.37	93	38672	3.946	ng	# 43
10) Phenol	6.37	94	46555	5.366	ng	# 50
17) 2-Methylphenol	6.97	107	47088m	8.717	ng	
20) 3+4-Methylphenols	7.13	107	109927	16.855	ng	# 83
22) Acetophenone	7.12	105	269752	32.520	ng	# 97
27) 2,4-Dimethylphenol	7.66	122	88181	18.491	ng	98
31) Naphthalene	8.03	128	5737516	342.213	ng	95
37) 2-Methylnaphthalene	8.70	142	479580	44.372	ng	98
45) 1,1'-Biphenyl	9.16	154	70399	4.468	ng	97
49) Dimethylphthalate	9.46	163	75327	5.355	ng	100
51) Acenaphthene	9.76	154	98203	8.553	ng	99
54) Dibenzofuran	9.94	168	132497	7.760	ng	98
57) Fluorene	10.28	166	57336	4.706	ng	99
70) Phenanthrene	11.23	178	116885	6.212	ng	99
72) Carbazole	11.44	167	67769	3.569	ng	100

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

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