

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100218\
 Data File : BF109495.D
 Acq On : 2 Oct 2018 12:35
 Operator : JU/SJ
 Sample : J5155-03DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW26S-GWDL

Manual Integrations
 APPROVED

Sohil
 10/3/2018 9:30:58 AM

Quant Time: Oct 02 14:09:10 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	102610	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	333672	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	193903	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	351222	20.00	ng	0.00
75) Chrysene-d12	13.84	240	262554	20.00	ng	0.00
86) Perylene-d12	15.24	264	270110	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	32432	5.21	ng	0.01
7) Phenol-d6	6.36	99	45356	5.71	ng	0.00
23) Nitrobenzene-d5	7.26	82	104922	18.56	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	42258	22.11	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	208392	16.21	ng	-0.02
78) Terphenyl-d14	12.79	244	157408	14.71	ng	-0.01
Target Compounds						
6) Aniline	6.37	93	58984	5.799	ng	Qvalue # 82
10) Phenol	6.37	94	32747m	3.638	ng	
17) 2-Methylphenol	6.97	107	47088m	8.400	ng	95
20) 3+4-Methylphenols	7.13	107	149930	22.154	ng	
27) 2,4-Dimethylphenol	7.66	122	300422	67.738	ng	99
31) Naphthalene	8.05	128	8046795	516.072	ng	94
37) 2-Methylnaphthalene	8.70	142	433498	43.127	ng	97
45) 1,1'-Biphenyl	9.16	154	164447	10.409	ng	98
49) Dimethylphthalate	9.46	163	53054	3.762	ng	# 99
51) Acenaphthene	9.77	154	295159	25.640	ng	
54) Dibenzofuran	9.94	168	334407	19.534	ng	99
57) Fluorene	10.28	166	205575	16.830	ng	100
70) Phenanthrene	11.24	178	314534	17.936	ng	99
71) Anthracene	11.29	178	42163	2.370	ng	98
72) Carbazole	11.44	167	227947	12.881	ng	100

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

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