

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100118\
 Data File : BF109488.D
 Acq On : 1 Oct 2018 22:09
 Operator : JU/SJ
 Sample : J5168-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW27S-GW

Manual Integrations
 APPROVED

Sohil
 10/2/2018 1:01:37 PM

Quant Time: Oct 02 10:20:11 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	96460	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	353722	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	149873	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	248638	20.00	ng	0.00
75) Chrysene-d12	13.84	240	233488	20.00	ng	0.00
86) Perylene-d12	15.24	264	172086	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	130732	22.32	ng	0.00
7) Phenol-d6	6.35	99	104953	14.04	ng	0.00
23) Nitrobenzene-d5	7.26	82	279524	46.65	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	73386	49.67	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	507475	51.07	ng	-0.01
78) Terphenyl-d14	12.79	244	400437	42.09	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.37	88	9021	3.345	ng	# 81
10) Phenol	6.36	94	23546	2.782	ng	83
27) 2,4-Dimethylphenol	7.65	122	58445	12.431	ng	96
31) Naphthalene	8.01	128	1187934	71.868	ng	99
37) 2-Methylnaphthalene	8.80	142	420829m	39.493	ng	
45) 1,1'-Biphenyl	9.16	154	59836	4.900	ng	97
48) Acenaphthylene	9.59	152	41067	2.791	ng	# 91
49) Dimethylphthalate	9.46	163	231920	21.276	ng	100
51) Acenaphthene	9.77	154	186932	21.009	ng	99
54) Dibenzofuran	9.94	168	179456	13.562	ng	98
57) Fluorene	10.28	166	121800	12.901	ng	98
70) Phenanthrene	11.24	178	168226	13.551	ng	98
71) Anthracene	11.29	178	32633	2.592	ng	98
72) Carbazole	11.45	167	221540	17.684	ng	99

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

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