

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100518\
 Data File : BF109619.D
 Acq On : 5 Oct 2018 20:07
 Operator : JU/SJ
 Sample : J5205-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-100418

Manual Integrations
 APPROVED

Sohil
 10/8/2018 12:21:23 PM

Quant Time: Oct 06 06:37:33 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	99000	20.00	ng	-0.01
21) Naphthalene-d8	7.98	136	401723	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	182135	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	297890	20.00	ng	0.00
75) Chrysene-d12	13.83	240	220889	20.00	ng	-0.01
86) Perylene-d12	15.24	264	223882	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	485794	80.82	ng	0.01
7) Phenol-d6	6.35	99	670092	87.37	ng	0.00
23) Nitrobenzene-d5	7.26	82	429978	63.18	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	149652	83.35	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	794738	65.81	ng	-0.02
78) Terphenyl-d14	12.79	244	555001	61.66	ng	-0.02
Target Compounds						
10) Phenol	6.36	94	18257	2.102	ng	Qvalue # 47
49) Dimethylphthalate	9.45	163	77107	5.821	ng	99
70) Phenanthrene	11.23	178	160579	10.796	ng	99
71) Anthracene	11.29	178	40907	2.712	ng	98
74) Fluoranthene	12.42	202	259785	16.344	ng	95
77) Pyrene	12.64	202	225848	14.266	ng	99
80) Benzo(a)anthracene	13.83	228	101277	7.612	ng	98
82) Chrysene	13.86	228	101169	7.672	ng	97
85) Indeno(1,2,3-cd)pyrene	16.59	276	41508	3.883	ng	# 87
87) Benzo(b)fluoranthene	14.84	252	115293m	9.372	ng	
88) Benzo(k)fluoranthene	14.86	252	49571m	4.246	ng	
89) Benzo(a)pyrene	15.18	252	87279	7.873	ng	98
91) Benzo(a,h,i)perylene	16.99	276	40317m	4.511	ng	

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

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