

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109403.D
 Acq On : 27 Sep 2018 22:21
 Operator : JU/SJ
 Sample : J5096-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MT-T-5-2

Manual Integrations
 APPROVED

Sohil
 9/28/2018 2:40:29 PM

Quant Time: Sep 28 05:48:21 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	6.71	152	119779	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	445028	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	187202	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	302385	20.00	ng	0.00
75) Chrysene-d12	13.84	240	280460	20.00	ng	0.00
86) Perylene-d12	15.24	264	203722	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	645062	88.70	ng	0.00
7) Phenol-d6	6.35	99	817830	88.13	ng	0.00
23) Nitrobenzene-d5	7.27	82	528423	70.09	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	167621	90.83	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	860574	69.33	ng	-0.01
78) Terphenyl-d14	12.80	244	621412	54.38	ng	0.00
Target Compounds						
10) Phenol	6.36	94	23193	2.207	ng	Qvalue # 51
49) Dimethylphthalate	9.46	163	277046	20.347	ng	100
70) Phenanthrene	11.24	178	59062	3.912	ng	97
74) Fluoranthene	12.42	202	124352	7.707	ng	93
77) Pyrene	12.64	202	101961	5.073	ng	99
80) Benzo(a)anthracene	13.83	228	51632	3.056	ng	98
82) Chrysene	13.86	228	59408	3.548	ng	96
87) Benzo(b)fluoranthene	14.84	252	59961m	5.356	ng	
89) Benzo(a)pyrene	15.18	252	33160	3.287	ng	# 97
91) Benzo(g,h,i)perylene	16.99	276	18718	2.301	ng	# 91

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

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