

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092018\
 Data File : BF109193.D
 Acq On : 20 Sep 2018 20:26
 Operator : JU/SJ
 Sample : J4987-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR200036

Manual Integrations
 APPROVED

Sohil
 9/21/2018 10:37:58 AM

Quant Time: Sep 21 07:45:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 14 13:26:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	89457	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	389778	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	222148	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	319038	20.00	ng	0.00
75) Chrysene-d12	13.85	240	201638	20.00	ng	0.00
86) Perylene-d12	15.26	264	228147	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	426351	79.16	ng	0.01
7) Phenol-d6	6.36	99	554331	79.82	ng	0.00
23) Nitrobenzene-d5	7.28	82	404944	58.27	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	150327	62.35	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	478226	33.73	ng	0.00
78) Terphenyl-d14	12.81	244	411103	48.33	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.47	163	237070	14.931	ng	99
70) Phenanthrene	11.25	178	76321	4.880	ng	99
74) Fluoranthene	12.44	202	403769	24.455	ng	96
77) Pyrene	12.67	202	557808	37.243	ng	99
80) Benzo(a)anthracene	13.84	228	312792	25.622	ng	98
82) Chrysene	13.88	228	352711	28.878	ng	99
85) Indeno(1,2,3-cd)pyrene	16.60	276	90539	9.272	ng	# 92
87) Benzo(b)fluoranthene	14.87	252	325531m	25.792	ng	
88) Benzo(k)fluoranthene	14.89	252	128036m	10.861	ng	
89) Benzo(a)pyrene	15.20	252	165946	14.805	ng	98
90) Dibenzo(a,h)anthracene	16.61	278	24935	2.648	ng	94
91) Benzo(a,h,i)perylene	17.01	276	66404	7.053	ng	94

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

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