

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF011218\
 Data File : BF102054.D
 Acq On : 12 Jan 2018 20:23
 Operator : SJ/JU
 Sample : J1094-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DBSW-2(0-4)

Manual Integrations
 APPROVED

Sohil
 1/15/2018 4:23:01 PM

Quant Time: Jan 12 23:50:26 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 11 12:53:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	223740	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	844905	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	324927	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	440284	20.00	ng	0.00
75) Chrysene-d12	13.89	240	360787	20.00	ng	0.00
86) Perylene-d12	15.31	264	295648	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1422175	89.76	ng	0.02
7) Phenol-d6	6.37	99	1663480	87.99	ng	0.00
23) Nitrobenzene-d5	7.30	82	1007624	70.08	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	218402	77.02	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	1359125	65.35	ng	0.00
78) Terphenyl-d14	12.83	244	787057	45.29	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.38	94	64669	3.027	ng	97
49) Dimethylphthalate	9.48	163	294784	11.309	ng	100
70) Phenanthrene	11.27	178	230726	8.971	ng	99
71) Anthracene	11.32	178	56673	2.173	ng	98
74) Fluoranthene	12.46	202	325141	12.850	ng	99
77) Pyrene	12.69	202	281847	8.838	ng	100
80) Benzo(a)anthracene	13.87	228	146429	6.361	ng	96
82) Chrysene	13.91	228	146056	6.085	ng	97
85) Indeno(1,2,3-cd)pyrene	16.69	276	53222	3.046	ng	96
87) Benzo(b)fluoranthene	14.90	252	160317m	8.316	ng	
88) Benzo(k)fluoranthene	14.92	252	59947m	3.224	ng	
89) Benzo(a)pyrene	15.24	252	106032	6.112	ng	# 87
91) Benzo(g,h,i)perylene	17.10	276	51995	3.729	ng	# 89

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

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