

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF021518\
 Data File : BF103011.D
 Acq On : 15 Feb 2018 12:15
 Operator : SJ/JU
 Sample : J1539-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MAIN-ST-LINE-1-WW@4

Manual Integrations
 APPROVED

Sohil
 2/16/2018 2:31:08 PM

Quant Time: Feb 15 14:22:09 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 15 14:08:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	198998	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	832580	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	377980	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	667224	20.00	ng	0.00
75) Chrysene-d12	14.04	240	525139	20.00	ng	0.00
86) Perylene-d12	15.53	264	449728	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1609617	122.84	ng	0.01
7) Phenol-d6	6.52	99	1920777	121.74	ng	0.00
23) Nitrobenzene-d5	7.45	82	1255801	104.63	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	425084	139.72	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1813331	79.61	ng	0.00
78) Terphenyl-d14	12.99	244	1673592	75.46	ng	0.00
Target Compounds						
10) Phenol	6.52	94	63207	3.738	ng	Qvalue # 72
49) Dimethylphthalate	9.62	163	165972	5.632	ng	99
70) Phenanthrene	11.43	178	323298	8.700	ng	98
71) Anthracene	11.48	178	83748	2.216	ng	99
74) Fluoranthene	12.62	202	341197	9.126	ng	98
77) Pyrene	12.84	202	296425	7.198	ng	99
80) Benzo(a)anthracene	14.03	228	142066	4.302	ng	97
82) Chrysene	14.07	228	129500	3.890	ng	96
87) Benzo(b)fluoranthene	15.09	252	118031m	4.048	ng	
89) Benzo(a)pyrene	15.46	252	83053	3.164	ng	# 92

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

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