

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031418\
 Data File : BF103654.D
 Acq On : 14 Mar 2018 21:19
 Operator : SJ/JU
 Sample : J1911-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-13-COMP

Manual Integrations
 APPROVED

Sohil
 3/15/2018 6:57:45 PM

Quant Time: Mar 15 01:39:36 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 14 18:08:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	135070	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	580143	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	241689	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	373453	20.00	ng	0.00
75) Chrysene-d12	14.06	240	241172	20.00	ng	0.00
86) Perylene-d12	15.57	264	209581	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	783541	87.74	ng	0.00
7) Phenol-d6	6.50	99	967516	88.54	ng	0.00
23) Nitrobenzene-d5	7.43	82	678254	66.77	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	141861	69.34	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	905352	60.07	ng	0.00
78) Terphenyl-d14	12.97	244	610263	60.83	ng	0.00
Target Compounds						
73) Di-n-butylphthalate	11.93	149	278237	10.595	ng	99
74) Fluoranthene	12.62	202	117522	5.690	ng	98
77) Pyrene	12.85	202	119139	6.336	ng	100
80) Benzo(a)anthracene	14.04	228	52745	3.371	ng	97
82) Chrysene	14.09	228	56831	3.740	ng	99
85) Indeno(1,2,3-cd)pyrene	17.13	276	32731	2.721	ng	# 91
87) Benzo(b)fluoranthene	15.13	252	77543m	5.627	ng	
88) Benzo(k)fluoranthene	15.15	252	28995m	2.235	ng	
89) Benzo(a)pyrene	15.51	252	55362	4.499	ng	# 93
91) Benzo(g,h,i)perylene	17.60	276	34565	3.719	ng	# 91

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

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