

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103470.D
 Acq On : 6 Mar 2018 21:55
 Operator : SJ/JU
 Sample : J1722-15 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-11

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:17:26 PM

Quant Time: Mar 07 01:19:00 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 11:52:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	130092	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	515891	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	216420	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	306117	20.00	ng	0.00
75) Chrysene-d12	14.06	240	209710	20.00	ng	0.00
86) Perylene-d12	15.58	264	157347	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	493215	57.34	ng	0.00
7) Phenol-d6	6.51	99	616246	58.55	ng	0.00
23) Nitrobenzene-d5	7.44	82	369108	40.86	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	62178	33.94	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	538829	36.91	ng	0.00
78) Terphenyl-d14	12.99	244	364501	41.78	ng	0.00
Target Compounds						
70) Phenanthrene	11.43	178	97108	5.764	ng	98
74) Fluoranthene	12.63	202	220774	13.041	ng	99
77) Pyrene	12.86	202	219122	13.402	ng	99
80) Benzo(a)anthracene	14.05	228	88741	6.522	ng	97
82) Chrysene	14.09	228	79886	6.047	ng	97
83) Bis(2-ethylhexyl)phthalate	14.01	149	135321	11.608	ng	# 98
85) Indeno(1,2,3-cd)pyrene	17.13	276	29764	2.846	ng	96
87) Benzo(b)fluoranthene	15.13	252	78351m	7.573	ng	
88) Benzo(k)fluoranthene	15.16	252	25803m	2.649	ng	
89) Benzo(a)pyrene	15.52	252	53926	5.837	ng	# 83
91) Benzo(g,h,i)perylene	17.60	276	29653	4.250	ng	# 88

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

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