

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103449.D
 Acq On : 6 Mar 2018 12:17
 Operator : SJ/JU
 Sample : J1723-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-26

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:14:43 PM

Quant Time: Mar 06 13:05:56 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	118249	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	495747	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	218473	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	348035	20.00	ng	0.00
75) Chrysene-d12	14.06	240	220296	20.00	ng	0.00
86) Perylene-d12	15.59	264	216186	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	868776	111.12	ng	0.00
7) Phenol-d6	6.51	99	1060354	110.83	ng	0.00
23) Nitrobenzene-d5	7.44	82	671150	77.31	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	154595	83.60	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	866238	64.41	ng	0.00
78) Terphenyl-d14	12.99	244	587639	64.12	ng	0.00
Target Compounds						
70) Phenanthrene	11.43	178	232620	12.145	ng	98
71) Anthracene	11.49	178	68472	3.486	ng	98
74) Fluoranthene	12.63	202	353848	18.384	ng	100
77) Pyrene	12.86	202	338983	19.736	ng	100
80) Benzo(a)anthracene	14.06	228	164723	11.524	ng	97
82) Chrysene	14.09	228	155494	11.204	ng	99
85) Indeno(1,2,3-cd)pyrene	17.13	276	74119	6.746	ng	98
87) Benzo(b)fluoranthene	15.14	252	213662m	15.032	ng	
88) Benzo(k)fluoranthene	15.16	252	58168m	4.346	ng	
89) Benzo(a)pyrene	15.52	252	139255	10.971	ng	# 92
90) Dibenzo(a,h)anthracene	17.12	278	20709	2.111	ng	# 75
91) Benzo(a,h,i)perylene	17.60	276	72483	7.561	ng	# 91

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

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