

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103421.D
 Acq On : 5 Mar 2018 20:56
 Operator : SJ/JU
 Sample : J1723-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:23:22 PM

Quant Time: Mar 06 00:51:20 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 13:40:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	145204	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	582263	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	242610	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	366458	20.00	ng	0.00
75) Chrysene-d12	14.07	240	218239	20.00	ng	0.00
86) Perylene-d12	15.59	264	167340	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1003974	104.57	ng	0.00
7) Phenol-d6	6.52	99	1222914	104.10	ng	0.00
23) Nitrobenzene-d5	7.44	82	721221	70.74	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	159232	77.54	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	972239	65.26	ng	0.00
78) Terphenyl-d14	13.00	244	638751	70.36	ng	0.00
Target Compounds						
10) Phenol	6.53	94	34695	2.544	ng	Qvalue # 66
49) Dimethylphthalate	9.62	163	90418	4.646	ng	99
70) Phenanthrene	11.43	178	192725	9.556	ng	99
71) Anthracene	11.49	178	63383	3.065	ng	98
74) Fluoranthene	12.63	202	387196	19.106	ng	99
77) Pyrene	12.86	202	382870	22.502	ng	99
80) Benzo(a)anthracene	14.06	228	197665	13.959	ng	98
82) Chrysene	14.10	228	180058	13.096	ng	98
85) Indeno(1,2,3-cd)pyrene	17.13	276	73771	6.777	ng	97
87) Benzo(b)fluoranthene	15.14	252	199213m	18.106	ng	
88) Benzo(k)fluoranthene	15.16	252	56605m	5.463	ng	
89) Benzo(a)pyrene	15.52	252	131340	13.368	ng	# 92
90) Dibenzo(a,h)anthracene	17.13	278	20531	2.704	ng	# 63
91) Benzo(a,h,i)perylene	17.61	276	71637	9.655	ng	# 89

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

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