

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
 Data File : BF103332.D
 Acq On : 1 Mar 2018 13:15
 Operator : SJ/JU
 Sample : J1720-15
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-29

Manual Integrations
 APPROVED

Sohil
 3/2/2018 3:30:28 PM

Quant Time: Mar 01 14:22:36 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	168147	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	672978	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	262099	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	366028	20.00	ng	0.00
75) Chrysene-d12	14.06	240	273175	20.00	ng	0.00
86) Perylene-d12	15.57	264	225571	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1123277	101.04	ng	0.01
7) Phenol-d6	6.52	99	1348000	99.09	ng	0.00
23) Nitrobenzene-d5	7.44	82	904399	76.75	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	180954	81.57	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1219014	78.71	ng	0.00
78) Terphenyl-d14	12.99	244	781737	68.79	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	135979	6.467	ng	100
70) Phenanthrene	11.43	178	216431	10.744	ng	99
71) Anthracene	11.48	178	64181	3.107	ng	99
74) Fluoranthene	12.62	202	393192	19.424	ng	99
77) Pyrene	12.86	202	378889	17.790	ng	99
80) Benzo(a)anthracene	14.04	228	203277	11.469	ng	97
82) Chrysene	14.09	228	188008	10.924	ng	97
85) Indeno(1,2,3-cd)pyrene	17.10	276	79269	5.818	ng	98
87) Benzo(b)fluoranthene	15.12	252	225698m	15.218	ng	
88) Benzo(k)fluoranthene	15.15	252	86312m	6.180	ng	
89) Benzo(a)pyrene	15.50	252	153100	11.560	ng	# 92
90) Dibenzo(a,h)anthracene	17.10	278	21652	2.116	ng	# 70
91) Benzo(a,h,i)perylene	17.57	276	78719	7.870	ng	# 91

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

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