

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030218\
 Data File : BF103375.D
 Acq On : 2 Mar 2018 19:44
 Operator : SJ/JU
 Sample : J1720-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-36

Manual Integrations
 APPROVED

Sohil
 3/5/2018 3:13:32 PM

Quant Time: Mar 02 23:01:08 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 02 13:28:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	141086	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	597485	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	247403	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	360655	20.00	ng	0.00
75) Chrysene-d12	14.06	240	232463	20.00	ng	0.00
86) Perylene-d12	15.57	264	203356	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	990582	106.19	ng	0.00
7) Phenol-d6	6.52	99	1178162	103.21	ng	0.00
23) Nitrobenzene-d5	7.44	82	791500	75.65	ng	-0.01
41) 2,4,6-Tribromophenol	10.71	330	183117	87.44	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1053846	70.41	ng	0.00
78) Terphenyl-d14	12.99	244	665369	68.80	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.63	163	122853	6.190	ng	99
70) Phenanthrene	11.43	178	282777	14.247	ng	98
71) Anthracene	11.49	178	83908	4.123	ng	100
74) Fluoranthene	12.63	202	463167	23.222	ng	98
77) Pyrene	12.86	202	432356	23.855	ng	100
80) Benzo(a)anthracene	14.05	228	222717	14.766	ng	98
82) Chrysene	14.09	228	211709	14.456	ng	98
85) Indeno(1,2,3-cd)pyrene	17.12	276	81481	7.028	ng	95
87) Benzo(b)fluoranthene	15.13	252	244793m	18.308	ng	
88) Benzo(k)fluoranthene	15.16	252	81220m	6.451	ng	
89) Benzo(a)pyrene	15.52	252	165590	13.869	ng	# 92
90) Dibenzo(a,h)anthracene	17.12	278	21748	2.357	ng	# 60
91) Benzo(a,h,i)perylene	17.59	276	77061	8.546	ng	# 89

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

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