

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103422.D
 Acq On : 5 Mar 2018 21:23
 Operator : SJ/JU
 Sample : J1724-17
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-32

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:23:26 PM

Quant Time: Mar 06 00:54:34 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	139371	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	563459	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	229834	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	342898	20.00	ng	0.00
75) Chrysene-d12	14.07	240	210897	20.00	ng	0.00
86) Perylene-d12	15.59	264	159800	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1034463	112.26	ng	0.00
7) Phenol-d6	6.52	99	1222945	108.46	ng	0.00
23) Nitrobenzene-d5	7.44	82	752016	76.22	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	159978	82.23	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1010338	73.25	ng	0.00
78) Terphenyl-d14	13.00	244	675942	77.05	ng	0.00
Target Compounds						
10) Phenol	6.53	94	50475	3.856	ng	Qvalue # 68
49) Dimethylphthalate	9.62	163	90907	4.931	ng	99
70) Phenanthrene	11.43	178	292692	15.510	ng	99
71) Anthracene	11.49	178	82766	4.277	ng	99
74) Fluoranthene	12.63	202	450387	23.751	ng	100
77) Pyrene	12.86	202	450112	27.374	ng	99
80) Benzo(a)anthracene	14.06	228	200607	14.660	ng	99
82) Chrysene	14.10	228	189946	14.296	ng	97
85) Indeno(1,2,3-cd)pyrene	17.14	276	66853	6.356	ng	97
87) Benzo(b)fluoranthene	15.14	252	201674m	19.195	ng	
88) Benzo(k)fluoranthene	15.17	252	57917m	5.854	ng	
89) Benzo(a)pyrene	15.53	252	127762	13.617	ng	# 89
90) Dibenzo(a,h)anthracene	17.14	278	18696	2.579	ng	# 57
91) Benzo(g,h,i)perylene	17.62	276	64839	9.151	ng	# 92

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

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