

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
 Data File : BF103455.D
 Acq On : 6 Mar 2018 14:57
 Operator : SJ/JU
 Sample : J1792-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-6

Manual Integrations
 APPROVED

Sohil
 3/7/2018 12:15:03 PM

Quant Time: Mar 06 16:39:27 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	114387	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	491814	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	202213	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	314598	20.00	ng	0.00
75) Chrysene-d12	14.07	240	195692	20.00	ng	0.00
86) Perylene-d12	15.58	264	169463	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	699460	92.48	ng	0.00
7) Phenol-d6	6.51	99	854741	92.36	ng	0.00
23) Nitrobenzene-d5	7.44	82	550296	63.90	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	126423	73.86	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	747005	59.06	ng	0.00
78) Terphenyl-d14	12.99	244	462809	56.85	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	11.43	178	55226	3.190	ng	98
73) Di-n-butylphthalate	11.95	149	216066	9.767	ng	98
74) Fluoranthene	12.63	202	75799	4.357	ng	99
77) Pyrene	12.86	202	65001	4.260	ng	99
80) Benzo(a)anthracene	14.06	228	32299	2.544	ng	# 92
82) Chrysene	14.09	228	31841	2.583	ng	# 96
87) Benzo(b)fluoranthene	15.13	252	36214m	3.250	ng	
89) Benzo(a)pyrene	15.52	252	25836	2.597	ng	# 68

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

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