

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111317\
 Data File : BF100534.D
 Acq On : 14 Nov 2017 9:26
 Operator : SJ/JU
 Sample : I6409-01DL 2X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-110917-01DL

Manual Integrations
 APPROVED

SOHIL
 11/14/2017 5:20:47 PM

Quant Time: Nov 14 11:06:23 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 14 02:37:14 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	137286	20.00	ng	-0.01
21) Naphthalene-d8	8.17	136	507250	20.00	ng	-0.02
38) Acenaphthene-d10	9.93	164	192359	20.00	ng	-0.01
63) Phenanthrene-d10	11.41	188	285773	20.00	ng	-0.01
75) Chrysene-d12	14.05	240	270262	20.00	ng	-0.02
86) Perylene-d12	15.53	264	210431	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	405803	47.38	ng	0.00
7) Phenol-d6	6.51	99	502226	47.55	ng	-0.01
23) Nitrobenzene-d5	7.45	82	294364	38.37	ng	-0.02
41) 2,4,6-Tribromophenol	10.71	330	82707	42.19	ng	-0.02
44) 2-Fluorobiphenyl	9.25	172	399738	31.64	ng	-0.02
78) Terphenyl-d14	12.99	244	231487	19.68	ng	-0.02
Target Compounds						
10) Phenol	6.52	94	24493	2.209	ng	Qvalue 97
31) Naphthalene	8.19	128	50581	2.070	ng	99
48) Acenaphthylene	9.79	152	65171	3.258	ng	99
49) Dimethylphthalate	9.63	163	37254	2.537	ng	# 98
57) Fluorene	10.47	166	26255	2.102	ng	# 94
70) Phenanthrene	11.44	178	378503	25.156	ng	99
71) Anthracene	11.49	178	137107	8.982	ng	100
74) Fluoranthene	12.63	202	917316	57.525	ng	98
77) Pyrene	12.86	202	847045	43.967	ng	100
80) Benzo(a)anthracene	14.04	228	565004	34.716	ng	96
82) Chrysene	14.08	228	488961	31.025	ng	98
85) Indeno(1,2,3-cd)pyrene	17.02	276	156383	12.268	ng	# 91
87) Benzo(b)fluoranthene	15.10	252	504930m	40.502	ng	
88) Benzo(k)fluoranthene	15.12	252	161530m	13.713	ng	
89) Benzo(a)pyrene	15.47	252	324679	29.376	ng	98
90) Dibenzo(a,h)anthracene	17.02	278	42468	4.698	ng	93
91) Benzo(a,h,i)perylene	17.47	276	138635	15.669	ng	# 93

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

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