

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111317\
 Data File : BF100508.D
 Acq On : 13 Nov 2017 18:13
 Operator : SJ/JU
 Sample : I6301-07 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-110917-D

Manual Integrations
 APPROVED

SOHIL
 11/14/2017 5:20:29 PM

Quant Time: Nov 14 02:40:16 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 13 14:08:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	84454	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	316586	20.00	ng	0.00
38) Acenaphthene-d10	9.93	164	126759	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	193688	20.00	ng	0.00
75) Chrysene-d12	14.06	240	182931	20.00	ng	0.00
86) Perylene-d12	15.53	264	139465	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	283075	53.73	ng	0.00
7) Phenol-d6	6.51	99	341686	52.59	ng	0.00
23) Nitrobenzene-d5	7.45	82	203372	42.47	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	69280	53.64	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	371830	44.67	ng	0.00
78) Terphenyl-d14	13.00	244	246189	30.92	ng	0.00
Target Compounds						
10) Phenol	6.52	94	14162	2.076	ng	88
49) Dimethylphthalate	9.64	163	61150	6.318	ng	# 99
70) Phenanthrene	11.44	178	113309	11.111	ng	98
71) Anthracene	11.49	178	34529	3.337	ng	96
74) Fluoranthene	12.63	202	190448	17.621	ng	95
77) Pyrene	12.86	202	175844	13.485	ng	99
80) Benzo(a)anthracene	14.04	228	96048m	8.719	ng	
82) Chrysene	14.08	228	88148	8.263	ng	95
85) Indeno(1,2,3-cd)pyrene	17.02	276	33459	3.878	ng	# 92
87) Benzo(b)fluoranthene	15.10	252	88246m	10.680	ng	
88) Benzo(k)fluoranthene	15.12	252	31098m	3.983	ng	
89) Benzo(a)pyrene	15.47	252	58899	8.041	ng	98
91) Benzo(a,h,i)perylene	17.47	276	36595	6.241	ng	92

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

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