

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112817\
 Data File : BF100936.D
 Acq On : 28 Nov 2017 16:19
 Operator : SJ/JU
 Sample : I6579-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HAR-4-E(7-8)

Manual Integrations
 APPROVED

Sohil
 11/29/2017 4:28:33 PM

Quant Time: Nov 29 06:49:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 15:21:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	59494	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	224757	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	97624	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	150532	20.00	ng	0.00
75) Chrysene-d12	14.02	240	136171	20.00	ng	0.00
86) Perylene-d12	15.50	264	143927	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	335813	90.48	ng	0.00
7) Phenol-d6	6.49	99	413474	90.34	ng	0.00
23) Nitrobenzene-d5	7.42	82	248729	73.16	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	90244	90.72	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	488338	76.17	ng	0.00
78) Terphenyl-d14	12.96	244	348937	58.88	ng	0.00
Target Compounds						
10) Phenol	6.50	94	9972	2.076	ng	90
48) Acenaphthylene	9.76	152	73427	7.233	ng	99
49) Dimethylphthalate	9.60	163	20533	2.755	ng	98
70) Phenanthrene	11.40	178	48939	6.175	ng	98
71) Anthracene	11.46	178	23680	2.945	ng	97
74) Fluoranthene	12.59	202	85945	10.232	ng	94
77) Pyrene	12.82	202	82842	8.534	ng	98
80) Benzo(a)anthracene	14.01	228	57725m	7.039	ng	
82) Chrysene	14.04	228	51755	6.518	ng	98
85) Indeno(1,2,3-cd)pyrene	16.97	276	61154	9.521	ng	# 90
87) Benzo(b)fluoranthene	15.06	252	102533m	12.025	ng	
88) Benzo(k)fluoranthene	15.09	252	33065m	4.104	ng	
89) Benzo(a)pyrene	15.43	252	77945	10.311	ng	95
90) Dibenzo(a,h)anthracene	16.97	278	15398m	2.491	ng	
91) Benzo(a,h,i)perylene	17.42	276	95939	15.853	ng	# 94

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

