

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110917\
 Data File : BF100356.D
 Acq On : 9 Nov 2017 22:23
 Operator : SJ/JU
 Sample : I6379-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-(2-8)

Manual Integrations
 APPROVED

Sohil
 11/10/2017 1:15:19 PM

Quant Time: Nov 10 02:47:09 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 16:04:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	192399	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	771088	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	356861	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	656309	20.00	ng	0.00
75) Chrysene-d12	14.06	240	406572	20.00	ng	0.00
86) Perylene-d12	15.54	264	409873	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1170293	97.50	ng	0.02
7) Phenol-d6	6.53	99	1468767	99.24	ng	0.00
23) Nitrobenzene-d5	7.46	82	940620	80.65	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	381989	105.05	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1668201	71.18	ng	0.00
78) Terphenyl-d14	13.01	244	1389696	78.54	ng	0.00
Target Compounds						
10) Phenol	6.53	94	51159	3.293	ng	Qvalue # 59
70) Phenanthrene	11.45	178	187934	5.439	ng	98
74) Fluoranthene	12.63	202	364843	9.962	ng	98
77) Pyrene	12.86	202	329914	11.383	ng	99
80) Benzo(a)anthracene	14.05	228	154473	6.309	ng	94
82) Chrysene	14.09	228	151904	6.407	ng	98
85) Indeno(1,2,3-cd)pyrene	17.03	276	103666	5.406	ng	# 91
87) Benzo(b)fluoranthene	15.11	252	192055m	7.909	ng	
88) Benzo(k)fluoranthene	15.13	252	78015m	3.400	ng	
89) Benzo(a)pyrene	15.48	252	144010	6.690	ng	98
91) Benzo(g,h,i)perylene	17.48	276	98788	5.732	ng	95

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

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