

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111417\
 Data File : BF100544.D
 Acq On : 14 Nov 2017 15:58
 Operator : SJ/JU
 Sample : I6301-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

SOHIL
 11/15/2017 12:30:06 PM

Quant Time: Nov 14 16:33:13 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 14:17:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	134294	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	495554	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	195715	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	312255	20.00	ng	0.00
75) Chrysene-d12	14.04	240	267568	20.00	ng	0.00
86) Perylene-d12	15.52	264	192967	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	780931	93.22	ng	0.01
7) Phenol-d6	6.51	99	955565	92.50	ng	0.00
23) Nitrobenzene-d5	7.45	82	580805	77.49	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	173415	86.95	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	887576	69.06	ng	0.00
78) Terphenyl-d14	12.99	244	626069	53.76	ng	0.00
Target Compounds						
10) Phenol	6.52	94	41756	3.850	ng	96
49) Dimethylphthalate	9.63	163	168000	11.243	ng	98
74) Fluoranthene	12.62	202	81117	4.655	ng	98
77) Pyrene	12.85	202	75229	3.944	ng	99
80) Benzo(a)anthracene	14.03	228	47908	2.973	ng	93
82) Chrysene	14.07	228	44497	2.852	ng	96
87) Benzo(b)fluoranthene	15.09	252	46470m	4.065	ng	
89) Benzo(a)pyrene	15.46	252	32012	3.159	ng	# 89
91) Benzo(g,h,i)perylene	17.45	276	17345	2.138	ng	95

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

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