

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100083.D
 Acq On : 2 Nov 2017 14:33
 Operator : SJ/JU
 Sample : I6236-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 133

Manual Integrations
 APPROVED

Sohil
 11/3/2017 11:19:16 AM

Quant Time: Nov 02 16:30:22 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 02 16:13:54 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	87909	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	362218	20.00	ng	0.00
38) Acenaphthene-d10	9.81	164	162548	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	260637	20.00	ng	0.00
75) Chrysene-d12	13.98	240	169635	20.00	ng	0.00
86) Perylene-d12	15.52	264	163989	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	268996	48.04	ng	0.00
7) Phenol-d6	6.42	99	602020	90.67	ng	0.00
23) Nitrobenzene-d5	7.34	82	399421	70.99	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	24023	16.22	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	810762	76.95	ng	0.00
78) Terphenyl-d14	12.89	244	595311	76.69	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.53	163	122063	9.482	ng	100
59) Diethylphthalate	10.22	149	35392	2.815	ng	97
70) Phenanthrene	11.33	178	153276	10.421	ng	97
71) Anthracene	11.38	178	46936	3.144	ng	96
74) Fluoranthene	12.52	202	178302	11.664	ng	100
77) Pyrene	12.76	202	145840	11.306	ng	98
80) Benzo(a)anthracene	13.97	228	66191	6.155	ng	97
82) Chrysene	14.01	228	54445	5.385	ng	97
85) Indeno(1,2,3-cd)pyrene	17.03	276	31861	3.433	ng	99
87) Benzo(b)fluoranthene	15.09	252	74802m	7.224	ng	
88) Benzo(k)fluoranthene	15.12	252	24624m	2.522	ng	
89) Benzo(a)pyrene	15.46	252	53678	5.771	ng	# 92
91) Benzo(a,h,i)perylene	17.49	276	32617	4.034	ng	# 90

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

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