

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110317\
 Data File : BF100166.D
 Acq On : 4 Nov 2017 6:41
 Operator : SJ/JU
 Sample : I6265-02
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:19:51 PM

Quant Time: Nov 06 09:17:56 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 15:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	91470	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	362900	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	153723	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	234798	20.00	ng	0.00
75) Chrysene-d12	13.95	240	151721	20.00	ng	0.00
86) Perylene-d12	15.42	264	152537	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	553139	94.94	ng	0.00
7) Phenol-d6	6.42	99	652367	94.43	ng	-0.01
23) Nitrobenzene-d5	7.34	82	356312	63.21	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	121340	86.62	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	601388	60.36	ng	0.00
78) Terphenyl-d14	12.87	244	361451	52.06	ng	0.00
Target Compounds						
10) Phenol	6.43	94	36003	4.747	ng	# 79
20) 3+4-Methylphenols	7.21	107	79529	13.150	ng	# 80
31) Naphthalene	8.07	128	193411	11.041	ng	97
49) Dimethylphthalate	9.52	163	60566	4.975	ng	99
70) Phenanthrene	11.32	178	159551	12.041	ng	98
71) Anthracene	11.38	178	36484	2.713	ng	99
74) Fluoranthene	12.52	202	162822	11.824	ng	97
77) Pyrene	12.74	202	149083	12.922	ng	98
80) Benzo(a)anthracene	13.94	228	60124	6.251	ng	95
82) Chrysene	13.98	228	56334	6.230	ng	97
85) Indeno(1,2,3-cd)pyrene	16.90	276	33073	3.984	ng	94
87) Benzo(b)fluoranthene	14.99	252	75567m	7.845	ng	
88) Benzo(k)fluoranthene	15.02	252	21463m	2.363	ng	
89) Benzo(a)pyrene	15.36	252	62281	7.198	ng	# 95
91) Benzo(a,h,i)perylene	17.35	276	29950	3.982	ng	97

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

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