

Data Path : Z:\HPCHEM1\BNA F\DATA\BF103017\
 Data File : BF099965.D
 Acq On : 30 Oct 2017 17:45
 Operator : SJ/JU
 Sample : I6171-05 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-102617-C

Manual Integrations
 APPROVED

Sohil
 10/31/2017 7:41:06 PM

Quant Time: Oct 31 02:51:43 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 15:58:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	97769	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	402265	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	182539	20.00	ng	-0.02
63) Phenanthrene-d10	11.32	188	308635	20.00	ng	-0.02
75) Chrysene-d12	13.97	240	193780	20.00	ng	-0.02
86) Perylene-d12	15.46	264	185315	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	320949	51.54	ng	-0.01
7) Phenol-d6	6.43	99	413433	55.99	ng	-0.01
23) Nitrobenzene-d5	7.36	82	213180	34.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	97657	58.71	ng	-0.02
44) 2-Fluorobiphenyl	9.15	172	510845	43.18	ng	-0.01
78) Terphenyl-d14	12.90	244	413674	46.65	ng	-0.02
Target Compounds						
10) Phenol	6.45	94	20188	2.490	ng	79
49) Dimethylphthalate	9.54	163	86701	5.998	ng	99
74) Fluoranthene	12.54	202	57439	3.173	ng	98
77) Pyrene	12.77	202	63148	4.286	ng	98
80) Benzo(a)anthracene	13.97	228	25297	2.059	ng	96
82) Chrysene	14.00	228	25864	2.240	ng	96
87) Benzo(b)fluoranthene	15.03	252	36572m	3.125	ng	
89) Benzo(a)pyrene	15.40	252	27618	2.627	ng	# 83
91) Benzo(g,h,i)perylene	17.42	276	20094	2.199	ng	# 85

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

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