

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
 Data File : BF099927.D
 Acq On : 28 Oct 2017 19:50
 Operator : SJ/JU
 Sample : I6166-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-72-11966

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:20:34 PM

Quant Time: Oct 30 04:16:54 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	109456	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	437130	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	184703	20.00	ng	-0.02
63) Phenanthrene-d10	11.32	188	263099	20.00	ng	-0.02
75) Chrysene-d12	13.97	240	182041	20.00	ng	-0.02
86) Perylene-d12	15.43	264	185455	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	459069	65.85	ng	0.00
7) Phenol-d6	6.43	99	542659	65.64	ng	-0.01
23) Nitrobenzene-d5	7.36	82	313949	46.24	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	99877	59.34	ng	-0.02
44) 2-Fluorobiphenyl	9.14	172	656499	54.84	ng	-0.02
78) Terphenyl-d14	12.90	244	387389	46.51	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.54	163	104334	7.133	ng	99
70) Phenanthrene	11.34	178	75671	5.096	ng	99
74) Fluoranthene	12.53	202	100513	6.514	ng	97
77) Pyrene	12.76	202	89758	6.484	ng	98
80) Benzo(a)anthracene	13.96	228	45238	3.920	ng	96
82) Chrysene	13.99	228	38467	3.546	ng	97
83) Bis(2-ethylhexyl)phthalate	13.92	149	53596	5.599	ng	98
85) Indeno(1,2,3-cd)pyrene	16.91	276	23789	2.388	ng	93
87) Benzo(b)fluoranthene	15.01	252	58806m	5.022	ng	
89) Benzo(a)pyrene	15.37	252	38870	3.695	ng	# 84
91) Benzo(g,h,i)perylene	17.36	276	22249	2.433	ng	# 81

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

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