

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122917\
 Data File : BF101819.D
 Acq On : 29 Dec 2017 10:46
 Operator : SJ/JU
 Sample : I7056-01 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-9(0-10)COMP

Manual Integrations
 APPROVED

Sohil
 1/2/2018 1:04:53 PM

Quant Time: Dec 29 14:33:40 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 29 13:21:45 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	155700	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	617319	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	258028	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	440300	20.00	ng	0.00
75) Chrysene-d12	13.97	240	286212	20.00	ng	0.00
86) Perylene-d12	15.43	264	303127	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	492732	47.14	ng	0.00
7) Phenol-d6	6.43	99	614489	47.24	ng	0.00
23) Nitrobenzene-d5	7.36	82	370312	33.39	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	115831	46.59	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	635972	38.23	ng	0.00
78) Terphenyl-d14	12.89	244	467336	39.36	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.54	163	66631	3.210	ng	98
70) Phenanthrene	11.34	178	242468	9.902	ng	99
71) Anthracene	11.39	178	76961	3.077	ng	97
74) Fluoranthene	12.53	202	616847	24.001	ng	98
77) Pyrene	12.76	202	526739	24.522	ng	98
80) Benzo(a)anthracene	13.96	228	312094	16.514	ng	99
82) Chrysene	13.99	228	260466	14.597	ng	100
85) Indeno(1,2,3-cd)pyrene	16.92	276	157109	9.887	ng	# 90
87) Benzo(b)fluoranthene	15.01	252	379037m	20.515	ng	
88) Benzo(k)fluoranthene	15.03	252	118197m	6.580	ng	
89) Benzo(a)pyrene	15.38	252	255561	15.004	ng	98
90) Dibenzo(a,h)anthracene	16.92	278	37474m	2.563	ng	
91) Benzo(a,h,i)perylene	17.37	276	139999	9.668	ng	# 94

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

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