

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120617\
 Data File : BF101191.D
 Acq On : 7 Dec 2017 2:45
 Operator : SJ/JU
 Sample : I6656-04 2X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BART-3-E(0-1)

Manual Integrations
 APPROVED

Sohil
 12/7/2017 11:36:18 AM

Quant Time: Dec 07 05:33:04 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 04 13:07:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	44883	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	162563	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	67714	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	107974	20.00	ng	0.00
75) Chrysene-d12	14.02	240	106039	20.00	ng	0.00
86) Perylene-d12	15.50	264	85757	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	119116	43.85	ng	0.00
7) Phenol-d6	6.47	99	144489	43.82	ng	0.00
23) Nitrobenzene-d5	7.41	82	87832	32.23	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	24671	30.33	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	192168	38.83	ng	-0.01
78) Terphenyl-d14	12.96	244	151151	31.18	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.60	163	14807	2.711	ng	Qvalue # 98
70) Phenanthrene	11.40	178	24304	4.087	ng	99
74) Fluoranthene	12.59	202	46293	7.024	ng	94
77) Pyrene	12.82	202	36950	5.050	ng	96
80) Benzo(a)anthracene	14.04	228	22827	3.409	ng	94
82) Chrysene	14.04	228	22827	3.671	ng	98
87) Benzo(b)fluoranthene	15.06	252	24862m	4.840	ng	
89) Benzo(a)pyrene	15.44	252	13277	2.843	ng	# 91
91) Benzo(g,h,i)perylene	17.44	276	8703	2.243	ng	# 89

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

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