

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
 Data File : BF103226.D
 Acq On : 26 Feb 2018 19:19
 Operator : SJ/JU
 Sample : J1691-09 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3

Manual Integrations
 APPROVED

Sohil
 2/27/2018 12:22:57 PM

Quant Time: Feb 27 00:39:18 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	174744	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	730673	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	317301	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	494177	20.00	ng	0.00
75) Chrysene-d12	14.04	240	309361	20.00	ng	0.00
86) Perylene-d12	15.54	264	308828	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	610304	52.82	ng	0.00
7) Phenol-d6	6.50	99	715069	50.58	ng	0.00
23) Nitrobenzene-d5	7.43	82	450425	35.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	126206	46.99	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	740550	34.22	ng	0.00
78) Terphenyl-d14	12.98	244	497185	38.63	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	52922	2.079	ng	Qvalue # 99
74) Fluoranthene	12.62	202	116246	4.254	ng	98
77) Pyrene	12.84	202	108906	4.515	ng	97
80) Benzo(a)anthracene	14.03	228	63288	3.153	ng	95
82) Chrysene	14.07	228	57036	2.926	ng	96
85) Indeno(1,2,3-cd)pyrene	17.05	276	33851	2.194	ng	97
87) Benzo(b)fluoranthene	15.10	252	89423m	4.404	ng	
89) Benzo(a)pyrene	15.47	252	56945	3.140	ng	# 87
91) Benzo(g,h,i)perylene	17.51	276	35849	2.618	ng	# 91

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

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