

Data Path : U:\HPCHEM1\BNA F\DATA\BF020518\
 Data File : BF102740.D
 Acq On : 5 Feb 2018 16:17
 Operator : SJ/JU
 Sample : J1444-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-PH3-MGP-03

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:25:32 PM

Quant Time: Feb 06 10:43:21 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 12:44:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	199782	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	816423	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	358400	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	544714	20.00	ng	0.00
75) Chrysene-d12	14.04	240	346369	20.00	ng	0.00
86) Perylene-d12	15.52	264	331741	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1185971	90.15	ng	0.00
7) Phenol-d6	6.50	99	1405976	88.76	ng	0.00
23) Nitrobenzene-d5	7.45	82	847001	71.97	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	246440	85.43	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1058363	49.00	ng	0.00
78) Terphenyl-d14	12.99	244	612111	41.84	ng	-0.01
Target Compounds						
10) Phenol	6.52	94	100977	5.949	ng	99
49) Dimethylphthalate	9.63	163	117170	4.193	ng	98
56) 2,4-Dinitrotoluene	10.12	165	1622	3.288	ng	# 70
71) Anthracene	11.48	178	86007	2.787	ng	99
74) Fluoranthene	12.62	202	520351	17.048	ng	98
77) Pyrene	12.85	202	644234	23.719	ng	99
80) Benzo(a)anthracene	14.03	228	174341m	8.003	ng	
82) Chrysene	14.07	228	144868m	6.597	ng	
85) Indeno(1,2,3-cd)pyrene	17.00	276	47103	2.586	ng	94
87) Benzo(b)fluoranthene	15.09	252	135710m	6.310	ng	
89) Benzo(a)pyrene	15.45	252	132897	6.865	ng	# 95
91) Benzo(a,h,i)perylene	17.45	276	53595	3.485	ng	96

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

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