

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031218\
 Data File : BF103575.D
 Acq On : 12 Mar 2018 12:13
 Operator : SJ/JU
 Sample : J1776-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-3FT

Manual Integrations
 APPROVED

Sohil
 3/13/2018 4:12:48 PM

Quant Time: Mar 12 13:50:24 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 09 11:47:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	149962	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	614321	20.00	ng	-0.01
38) Acenaphthene-d10	9.90	164	259176	20.00	ng	-0.01
63) Phenanthrene-d10	11.40	188	401579	20.00	ng	0.00
75) Chrysene-d12	14.06	240	232587	20.00	ng	0.00
86) Perylene-d12	15.57	264	238033	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1074565	108.37	ng	0.01
7) Phenol-d6	6.51	99	1333314	109.89	ng	0.00
23) Nitrobenzene-d5	7.43	82	829446	77.11	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	188571	85.96	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1167813	75.58	ng	0.00
78) Terphenyl-d14	12.98	244	755205	78.05	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.61	163	119675	5.756	ng	Qvalue 100
70) Phenanthrene	11.42	178	83695	3.787	ng	98
74) Fluoranthene	12.62	202	104492	4.705	ng	98
77) Pyrene	12.85	202	95577	5.271	ng	98
80) Benzo(a)anthracene	14.04	228	33996	2.253	ng	97
82) Chrysene	14.08	228	37856	2.583	ng	96
85) Indeno(1,2,3-cd)pyrene	17.13	276	27334	2.356	ng	93
87) Benzo(b)fluoranthene	15.13	252	50526m	3.228	ng	
89) Benzo(a)pyrene	15.51	252	39916	2.856	ng	# 86
91) Benzo(g,h,i)perylene	17.60	276	31727	3.006	ng	# 83

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

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