

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050916\
 Data File : BF086848.D
 Acq On : 10 May 2016 4:16
 Operator : UM/SJ
 Sample : H2895-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP15-3-5FT

Manual Integrations
 APPROVED

sohil
 5/10/2016 11:06:15 AM

Quant Time: May 10 05:43:59 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 09 16:04:56 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	156692	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	663295	20.00	ng	-0.01
38) Acenaphthene-d10	9.72	164	301200	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	561368	20.00	ng	0.00
75) Chrysene-d12	13.83	240	323873	20.00	ng	-0.01
86) Perylene-d12	15.21	264	340784	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1270121	137.27	ng	0.01
7) Phenol-d6	6.35	99	1564312	126.50	ng	-0.01
23) Nitrobenzene-d5	7.26	82	1058702	80.15	ng	0.00
41) 2,4,6-Tribromophenol	10.52	330	391146	122.67	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	1172909	65.45	ng	0.00
78) Terphenyl-d14	12.78	244	783024	51.30	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.46	163	240898	10.48	ng	98
74) Fluoranthene	12.41	202	147042	4.77	ng	99
77) Pyrene	12.64	202	141101	5.69	ng	98
80) Benzo(a)anthracene	13.81	228	69149m	3.80	ng	
82) Chrysene	13.86	228	59906	3.24	ng	95
85) Indeno(1,2,3-cd)pyrene	16.49	276	45877	2.98	ng	# 87
87) Benzo(b)fluoranthene	14.83	252	95745m	4.32	ng	
89) Benzo(a)pyrene	15.15	252	68244	3.57	ng	# 93
91) Benzo(g,h,i)perylene	16.88	276	41314	2.53	ng	96

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

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