

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\
 Data File : BF088433.D
 Acq On : 26 Jun 2016 22:20
 Operator : UM/SJ
 Sample : H3663-10 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4

Manual Integrations
 APPROVED

sohil
 6/27/2016 8:02:32 PM

Quant Time: Jun 27 18:49:55 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 14:55:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.55	152	78604	20.00	ng	-0.06
21) Naphthalene-d8	7.84	136	311952	20.00	ng	-0.06
38) Acenaphthene-d10	9.58	164	140406	20.00	ng	-0.07
63) Phenanthrene-d10	11.05	188	232711	20.00	ng	-0.07
75) Chrysene-d12	13.68	240	174162	20.00	ng	-0.07
86) Perylene-d12	15.03	264	153909	20.00	ng	-0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	209312	45.52	ng	-0.05
7) Phenol-d6	6.22	99	268605	48.20	ng	-0.05
23) Nitrobenzene-d5	7.12	82	168773	31.59	ng	-0.06
41) 2,4,6-Tribromophenol	10.36	330	54371	36.67	ng	-0.07
44) 2-Fluorobiphenyl	8.91	172	331966	33.81	ng	-0.06
78) Terphenyl-d14	12.63	244	226827	29.64	ng	-0.07
Target Compounds						Qvalue
49) Dimethylphthalate	9.31	163	133455	13.12	ng	100
70) Phenanthrene	11.07	178	132732	10.87	ng	98
71) Anthracene	11.13	178	35896	2.91	ng	97
74) Fluoranthene	12.26	202	201152	15.61	ng	93
77) Pyrene	12.48	202	177188	13.71	ng	98
80) Benzo(a)anthracene	13.67	228	118514	11.94	ng	99
82) Chrysene	13.70	228	84799	9.08	ng	97
85) Indeno(1,2,3-cd)pyrene	16.25	276	55679	6.42	ng	# 100
87) Benzo(b)fluoranthene	14.67	252	145302m	13.55	ng	
88) Benzo(k)fluoranthene	14.69	252	46773m	5.78	ng	
89) Benzo(a)pyrene	14.98	252	96941m	11.17	ng	
91) Benzo(a,h,i)perylene	16.62	276	51958	6.27	ng	95

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

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