

Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
 Data File : BF089477.D
 Acq On : 31 Jul 2016 18:19
 Operator : UM/SJ
 Sample : H4252-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RU001

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:17:36 PM

Quant Time: Aug 01 05:51:22 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 04:11:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.49	152	46723	20.00	ng	-0.01
21) Naphthalene-d8	7.78	136	171997	20.00	ng	-0.01
38) Acenaphthene-d10	9.52	164	75104	20.00	ng	-0.01
63) Phenanthrene-d10	10.99	188	113941	20.00	ng	-0.01
75) Chrysene-d12	13.61	240	89834	20.00	ng	-0.01
86) Perylene-d12	14.94	264	101943	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.06	112	309436	105.75	ng	0.00
7) Phenol-d6	6.16	99	412559	106.68	ng	-0.01
23) Nitrobenzene-d5	7.06	82	244888	84.04	ng	-0.02
41) 2,4,6-Tribromophenol	10.31	330	58928	94.74	ng	-0.01
44) 2-Fluorobiphenyl	8.86	172	409151	79.95	ng	-0.01
78) Terphenyl-d14	12.57	244	259076	67.50	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.26	163	79595	13.93	ng	99
74) Fluoranthene	12.19	202	18280	2.96	ng	93
77) Pyrene	12.42	202	17954	2.64	ng	94
80) Benzo(a)anthracene	13.60	228	12226	2.41	ng	96
85) Indeno(1,2,3-cd)pyrene	16.13	276	8733	2.28	ng	95
87) Benzo(b)fluoranthene	14.61	252	16546m	2.61	ng	
89) Benzo(a)pyrene	14.89	252	11708	2.06	ng	95

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

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