

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
 Data File : BF088391.D
 Acq On : 25 Jun 2016 22:03
 Operator : UM/SJ
 Sample : H3629-06
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOCATION-5D(9-11)

Quant Time: Jun 26 03:47:32 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	6.56	152	73188	20.00	ng	-0.05
21) Naphthalene-d8	7.85	136	277700	20.00	ng	-0.05
38) Acenaphthene-d10	9.59	164	120770	20.00	ng	-0.06
63) Phenanthrene-d10	11.06	188	191616	20.00	ng	-0.06
75) Chrysene-d12	13.69	240	154872	20.00	ng	-0.06
86) Perylene-d12	15.04	264	120318	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.14	112	524459	122.50	ng	-0.03
7) Phenol-d6	6.23	99	628294	121.10	ng	-0.03
23) Nitrobenzene-d5	7.14	82	410976	86.42	ng	-0.03
41) 2,4,6-Tribromophenol	10.39	330	129996	101.94	ng	-0.05
44) 2-Fluorobiphenyl	8.92	172	646307	84.00	ng	-0.05
78) Terphenyl-d14	12.65	244	453864	66.69	ng	-0.05
Target Compounds						Qvalue
49) Dimethylphthalate	9.32	163	114120	13.04	ng	99
70) Phenanthrene	11.08	178	39595	3.94	ng	99
74) Fluoranthene	12.27	202	53670	5.06	ng	95
77) Pyrene	12.49	202	56627	4.93	ng	96
80) Benzo(a)anthracene	13.68	228	26637	3.02	ng	98
82) Chrysene	13.71	228	21585	2.60	ng	98
87) Benzo(b)fluoranthene	14.69	252	20537m	2.45	ng	
89) Benzo(a)pyrene	14.99	252	17214	2.54	ng	98

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

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