

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\
 Data File : BF088482.D
 Acq On : 28 Jun 2016 7:31
 Operator : UM/SJ
 Sample : H3602-07DL 10X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q8-B4(0-2)CDL

Quant Time: Jun 28 15:02:01 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 28 05:25:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.54	152	77000	20.00	ng	-0.01
21) Naphthalene-d8	7.83	136	311285	20.00	ng	-0.01
38) Acenaphthene-d10	9.58	164	128821	20.00	ng	-0.01
63) Phenanthrene-d10	11.05	188	192510	20.00	ng	-0.01
75) Chrysene-d12	13.68	240	168638	20.00	ng	-0.01
86) Perylene-d12	15.04	264	161772	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	55910	12.41	ng	0.00
7) Phenol-d6	6.22	99	75745	13.88	ng	-0.01
23) Nitrobenzene-d5	7.12	82	45376	8.51	ng	-0.01
41) 2,4,6-Tribromophenol	10.36	330	9962	7.32	ng	-0.01
44) 2-Fluorobiphenyl	8.90	172	72646	3.55	ng	-0.02
78) Terphenyl-d14	12.63	244	45523	6.14	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	1.69	88	74043	33.83	ng	# 87
51) Acenaphthene	9.61	154	490907	59.35	ng	98
54) Dibenzofuran	9.78	168	414497	36.39	ng	92
55) 4-Nitrophenol	9.78	139	151702	79.26	ng	# 6
57) Fluorene	10.12	166	410464	54.32	ng	97
70) Phenanthrene	11.09	178	1077673	106.68	ng	99
71) Anthracene	11.13	178	313179	30.73	ng	99
72) Carbazole	11.30	167	66414	6.96	ng	98
74) Fluoranthene	12.26	202	673410	63.17	ng	99
77) Pyrene	12.49	202	495799	39.62	ng	99
80) Benzo(a)anthracene	13.67	228	114053m	11.87	ng	
82) Chrysene	13.71	228	79791m	8.83	ng	
87) Benzo(b)fluoranthene	14.69	252	53193m	4.72	ng	
88) Benzo(k)fluoranthene	14.71	252	19828m	2.33	ng	
89) Benzo(a)pyrene	14.99	252	28597	3.13	ng	97

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

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