

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
 Data File : BF088382.D
 Acq On : 25 Jun 2016 17:43
 Operator : UM/SJ
 Sample : H3624-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-STA-1476(0-4)

Quant Time: Jun 26 03:24:34 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	77079	20.00	ng	-0.05
21) Naphthalene-d8	7.85	136	300945	20.00	ng	-0.05
38) Acenaphthene-d10	9.59	164	129130	20.00	ng	-0.06
63) Phenanthrene-d10	11.06	188	217077	20.00	ng	-0.06
75) Chrysene-d12	13.69	240	156491	20.00	ng	-0.06
86) Perylene-d12	15.05	264	138017	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.14	112	519585	115.24	ng	-0.03
7) Phenol-d6	6.23	99	615034	112.56	ng	-0.03
23) Nitrobenzene-d5	7.14	82	411151	79.78	ng	-0.03
41) 2,4,6-Tribromophenol	10.39	330	137928	101.16	ng	-0.05
44) 2-Fluorobiphenyl	8.92	172	692307	84.17	ng	-0.05
78) Terphenyl-d14	12.65	244	506845	73.70	ng	-0.05
Target Compounds						Qvalue
48) Acenaphthylene	9.45	152	55147	4.15	ng	98
49) Dimethylphthalate	9.32	163	104639	11.19	ng	99
70) Phenanthrene	11.08	178	129437	11.36	ng	98
71) Anthracene	11.14	178	57363	4.99	ng	98
74) Fluoranthene	12.27	202	328117	27.29	ng	97
77) Pyrene	12.50	202	260239	22.41	ng	97
80) Benzo(a)anthracene	13.68	228	177869	19.95	ng	97
82) Chrysene	13.71	228	131749m	15.70	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	52118	6.69	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	158696m	16.50	ng	
88) Benzo(k)fluoranthene	14.70	252	76449m	10.54	ng	
89) Benzo(a)pyrene	14.99	252	102454	13.16	ng	98
90) Dibenzo(a,h)anthracene	16.29	278	17196	2.38	ng	92
91) Benzo(g,h,i)perylene	16.65	276	40279	5.42	ng	97

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

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