

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
 Data File : BF088385.D
 Acq On : 25 Jun 2016 19:10
 Operator : UM/SJ
 Sample : H3622-13
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-STA-1461(0-4)

Quant Time: Jun 26 03:33:52 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	75861	20.00	ng	-0.05
21) Naphthalene-d8	7.85	136	291382	20.00	ng	-0.05
38) Acenaphthene-d10	9.59	164	124659	20.00	ng	-0.06
63) Phenanthrene-d10	11.06	188	205305	20.00	ng	-0.06
75) Chrysene-d12	13.69	240	159162	20.00	ng	-0.06
86) Perylene-d12	15.04	264	131758	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.15	112	501973	113.12	ng	-0.02
7) Phenol-d6	6.23	99	629448	117.04	ng	-0.03
23) Nitrobenzene-d5	7.14	82	463880	92.96	ng	-0.03
41) 2,4,6-Tribromophenol	10.39	330	127290	96.71	ng	-0.05
44) 2-Fluorobiphenyl	8.92	172	666232	83.88	ng	-0.05
78) Terphenyl-d14	12.65	244	436592	62.42	ng	-0.05
Target Compounds						Qvalue
31) Naphthalene	7.87	128	61510	4.27	ng	99
37) 2-Methylnaphthalene	8.56	142	38952	4.10	ng	97
48) Acenaphthylene	9.45	152	33286	2.59	ng	98
49) Dimethylphthalate	9.32	163	153185	16.96	ng	99
70) Phenanthrene	11.08	178	73352	6.81	ng	99
71) Anthracene	11.14	178	25718	2.37	ng	98
74) Fluoranthene	12.27	202	177164	15.58	ng	96
77) Pyrene	12.49	202	148738	12.59	ng	98
80) Benzo(a)anthracene	13.68	228	105713	11.66	ng	95
82) Chrysene	13.71	228	92268m	10.81	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	37222	4.70	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	113628m	12.37	ng	
88) Benzo(k)fluoranthene	14.70	252	51615m	7.45	ng	
89) Benzo(a)pyrene	14.99	252	63546	8.55	ng	97
91) Benzo(a,h,i)perylene	16.65	276	30972	4.36	ng	97

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

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