

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\
 Data File : BF088480.D
 Acq On : 28 Jun 2016 6:33
 Operator : UM/SJ
 Sample : H3622-19DL 2X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-STA-1471(0-4)DL

Quant Time: Jun 28 07:37:42 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	72370m	20.00	ng	-0.01
21) Naphthalene-d8	7.83	136	292464	20.00	ng	-0.01
38) Acenaphthene-d10	9.58	164	120640	20.00	ng	-0.01
63) Phenanthrene-d10	11.05	188	183537	20.00	ng	-0.01
75) Chrysene-d12	13.69	240	144444	20.00	ng	0.00
86) Perylene-d12	15.04	264	162984	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	303389	71.67	ng	-0.01
7) Phenol-d6	6.20	99	381870	74.43	ng	-0.02
23) Nitrobenzene-d5	7.12	82	219745	43.87	ng	-0.01
41) 2,4,6-Tribromophenol	10.36	330	53193	41.76	ng	-0.01
44) 2-Fluorobiphenyl	8.90	172	364235	44.81	ng	-0.02
78) Terphenyl-d14	12.63	244	193103	30.42	ng	-0.01
Target Compounds						Qvalue
48) Acenaphthylene	9.44	152	86522	6.97	ng	98
49) Dimethylphthalate	9.31	163	64670	7.40	ng	98
70) Phenanthrene	11.07	178	96486	10.02	ng	98
71) Anthracene	11.13	178	90371	9.30	ng	98
74) Fluoranthene	12.26	202	509958	50.17	ng	99
77) Pyrene	12.49	202	448483	41.84	ng	99
80) Benzo(a)anthracene	13.68	228	408951	49.68	ng	95
82) Chrysene	13.71	228	240208m	31.02	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	88768	12.34	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	346673m	30.52	ng	
88) Benzo(k)fluoranthene	14.70	252	136296m	15.91	ng	
89) Benzo(a)pyrene	14.99	252	225555	24.54	ng	97
90) Dibenzo(a,h)anthracene	16.27	278	36612	4.29	ng	# 91
91) Benzo(g,h,i)perylene	16.65	276	63401	7.22	ng	95

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

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