

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
 Data File : BF080149.D
 Acq On : 25 Jun 2015 1:39
 Operator : TP/IZ
 Sample : G2747-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-76(8-10)

Quant Time: Jun 25 04:58:06 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 00:52:28 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.32	152	46265	20.00	ng	0.00
21) Naphthalene-d8	8.61	136	171445	20.00	ng	0.00
38) Acenaphthene-d10	10.38	164	86981	20.00	ng	0.00
63) Phenanthrene-d10	11.88	188	180577	20.00	ng	-0.01
75) Chrysene-d12	14.77	240	184799	20.00	ng	-0.16
86) Perylene-d12	16.60	264	190848	20.00	ng	-0.24
System Monitoring Compounds						
5) 2-Fluorophenol	5.92	112	291875	111.68	ng	0.00
7) Phenol-d6	6.92	99	403595	116.04	ng	0.00
23) Nitrobenzene-d5	7.86	82	213465	72.48	ng	-0.01
41) 2,4,6-Tribromophenol	11.17	330	100418	118.06	ng	0.00
44) 2-Fluorobiphenyl	9.68	172	469922	77.05	ng	0.00
78) Terphenyl-d14	13.53	244	539881	68.23	ng	-0.08
Target Compounds						Qvalue
48) Acenaphthylene	10.24	152	60514	6.31	ng	98
49) Dimethylphthalate	10.07	163	42922	6.56	ng	# 97
70) Phenanthrene	11.90	178	32822	3.00	ng	96
71) Anthracene	11.96	178	27509	2.61	ng	98
74) Fluoranthene	13.15	202	72463	6.22	ng	87
77) Pyrene	13.39	202	116661m	10.25	ng	
80) Benzo(a)anthracene	14.76	228	84328	7.49	ng	# 68
82) Chrysene	14.80	228	96228m	9.27	ng	
85) Indeno(1,2,3-cd)pyrene	18.43	276	92591	7.07	ng	# 87
87) Benzo(b)fluoranthene	16.07	252	117273m	10.15	ng	
88) Benzo(k)fluoranthene	16.09	252	48939m	4.16	ng	
89) Benzo(a)pyrene	16.52	252	103801	9.46	ng	# 87
90) Dibenzo(a,h)anthracene	18.44	278	40224	3.58	ng	# 80
91) Benzo(g,h,i)perylene	18.97	276	100818	8.89	ng	# 77

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

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