

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022415\
 Data File : BF077414.D
 Acq On : 24 Feb 2015 13:38
 Operator : IZ / TP
 Sample : G1345-08DL 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 UB10PX3DL

Manual Integrations
 APPROVED

apatel
 2/25/2015 3:33:18 PM

Quant Time: Feb 25 10:33:39 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 19 16:57:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.30	152	45434	20.00	ng	-0.03
21) Naphthalene-d8	8.88	136	186869	20.00	ng	-0.03
38) Acenaphthene-d10	11.05	164	95323	20.00	ng	-0.05
63) Phenanthrene-d10	12.90	188	199480	20.00	ng	-0.03
75) Chrysene-d12	16.21	240	210785	20.00	ng	-0.01
86) Perylene-d12	18.04	264	237039	20.00	ng	0.09
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	141227	51.89	ng	-0.03
7) Phenol-d6	6.85	99	186296	53.24	ng	-0.02
23) Nitrobenzene-d5	7.98	82	101398	33.74	ng	-0.05
41) 2,4,6-Tribromophenol	12.03	330	49229	53.64	ng	-0.03
44) 2-Fluorobiphenyl	10.23	172	246167	40.27	ng	-0.03
78) Terphenyl-d14	14.89	244	271929	30.16	ng	-0.05
Target Compounds						Qvalue
31) Naphthalene	8.90	128	21346	2.28	ng	98
48) Acenaphthylene	10.88	152	20415	2.28	ng	98
49) Dimethylphthalate	10.73	163	18842	2.81	ng	98
51) Acenaphthene	11.09	154	21299	3.98	ng	98
54) Dibenzofuran	11.31	168	27026	3.27	ng	95
57) Fluorene	11.73	166	35160	5.32	ng	99
70) Phenanthrene	12.92	178	497877	43.06	ng	99
71) Anthracene	12.99	178	112719	9.82	ng	98
72) Carbazole	13.20	167	63922	5.86	ng	99
74) Fluoranthene	14.41	202	645695	51.13	ng	98
77) Pyrene	14.68	202	530975	41.18	ng	97
80) Benzo(a)anthracene	16.20	228	302709	24.50	ng	100
82) Chrysene	16.24	228	272215	23.47	ng	97
85) Indeno(1,2,3-cd)pyrene	19.53	276	195263m	14.76	ng	
87) Benzo(b)fluoranthene	17.56	252	384508m	27.90	ng	
88) Benzo(k)fluoranthene	17.60	252	107506m	7.96	ng	
89) Benzo(a)pyrene	17.96	252	268159	20.43	ng	# 96
90) Dibenzo(a,h)anthracene	19.55	278	42739	3.41	ng	# 87
91) Benzo(g,h,i)perylene	19.96	276	182303	14.43	ng	97

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

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