

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070715\
 Data File : BF080372.D
 Acq On : 7 Jul 2015 12:45
 Operator : TP/UM
 Sample : G2874-01DL 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1DL

Manual Integrations
 APPROVED

apatel
 7/8/2015 4:26:46 PM

Quant Time: Jul 07 14:41:52 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 07 14:37:43 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.23	152	37346	20.00	ng	0.00
21) Naphthalene-d8	8.52	136	141156	20.00	ng	0.00
38) Acenaphthene-d10	10.29	164	64675	20.00	ng	0.00
63) Phenanthrene-d10	11.80	188	138358	20.00	ng	0.00
75) Chrysene-d12	14.87	240	151217	20.00	ng	-0.01
86) Perylene-d12	16.81	264	151296	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.84	112	120673	59.18	ng	0.00
7) Phenol-d6	6.83	99	155995	57.71	ng	-0.01
23) Nitrobenzene-d5	7.79	82	85280	35.24	ng	0.00
41) 2,4,6-Tribromophenol	11.08	330	39614	58.54	ng	0.00
44) 2-Fluorobiphenyl	9.60	172	216273	45.13	ng	0.00
78) Terphenyl-d14	13.54	244	244738	37.75	ng	-0.01
Target Compounds						Qvalue
48) Acenaphthylene	10.14	152	20362	2.79	ng	99
49) Dimethylphthalate	9.98	163	33084	6.37	ng	100
70) Phenanthrene	11.82	178	104983	12.65	ng	98
71) Anthracene	11.88	178	55799	6.86	ng	97
72) Carbazole	12.03	167	36488	4.81	ng	100
74) Fluoranthene	13.12	202	488604	54.80	ng	98
77) Pyrene	13.38	202	433827	47.43	ng	99
80) Benzo(a)anthracene	14.85	228	253085m	27.77	ng	
82) Chrysene	14.90	228	200328	23.49	ng	98
85) Indeno(1,2,3-cd)pyrene	18.59	276	166197	16.58	ng	96
87) Benzo(b)fluoranthene	16.28	252	321176m	33.78	ng	
88) Benzo(k)fluoranthene	16.32	252	78177m	8.69	ng	
89) Benzo(a)pyrene	16.73	252	211068	24.27	ng	98
90) Dibenzo(a,h)anthracene	18.61	278	29641	3.46	ng	95
91) Benzo(g,h,i)perylene	19.14	276	164917	18.99	ng	98

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

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