

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
 Data File : BG017533.D
 Acq On : 7 Jun 2015 13:59
 Operator : TP/IZ
 Sample : G2457-13 2X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP-7

Manual Integrations
 APPROVED

apatel
 6/8/2015 6:01:06 PM

Quant Time: Jun 08 08:15:59 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 04 01:31:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	26424	20.00	ng	0.00
21) Naphthalene-d8	10.42	136	103665	20.00	ng	0.00
38) Acenaphthene-d10	14.29	164	61033	20.00	ng	0.00
63) Phenanthrene-d10	17.04	188	121482	20.00	ng	0.00
75) Chrysene-d12	21.24	240	139273	20.00	ng	0.00
86) Perylene-d12	23.48	264	139271	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	106892	71.27	ng	0.00
7) Phenol-d6	6.84	99	140443	68.73	ng	0.00
23) Nitrobenzene-d5	8.79	82	102426	49.32	ng	0.00
41) 2,4,6-Tribromophenol	15.80	330	52935	62.44	ng	0.00
44) 2-Fluorobiphenyl	12.90	172	197493	48.81	ng	0.00
78) Terphenyl-d14	19.68	244	242103	35.93	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.75	163	25681	4.83	ng	99
57) Fluorene	15.34	166	11737	2.27	ng	# 90
70) Phenanthrene	17.09	178	143521	21.19	ng	97
71) Anthracene	17.18	178	35538	5.25	ng	95
74) Fluoranthene	19.11	202	189043	21.34	ng	99
77) Pyrene	19.47	202	169349	19.74	ng	98
80) Benzo(a)anthracene	21.22	228	90646	10.58	ng	97
82) Chrysene	21.27	228	76952	9.92	ng	97
85) Indeno(1,2,3-cd)pyrene	25.75	276	35479	3.64	ng	97
87) Benzo(b)fluoranthene	22.81	252	96916	10.98	ng	# 97
88) Benzo(k)fluoranthene	22.84	252	31571m	3.74	ng	
89) Benzo(a)pyrene	23.38	252	74465	9.20	ng	96
91) Benzo(g,h,i)perylene	26.46	276	32322	4.04	ng	96

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

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