

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090815\
 Data File : BF081535.D
 Acq On : 8 Sep 2015 21:54
 Operator : UM/IZ
 Sample : G3555-12DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-2DL

Manual Integrations
 APPROVED

MMDadoda
 9/9/2015 2:59:06 PM

Quant Time: Sep 09 02:56:08 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 23:43:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.34	152	46246	20.00	ng	0.00
21) Naphthalene-d8	7.63	136	186729	20.00	ng	-0.01
38) Acenaphthene-d10	9.37	164	80527	20.00	ng	-0.01
63) Phenanthrene-d10	10.83	188	143072	20.00	ng	-0.01
75) Chrysene-d12	13.44	240	98007	20.00	ng	-0.01
86) Perylene-d12	14.75	264	93289	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.88	112	77024	25.84	ng	0.00
7) Phenol-d6	6.01	99	109914	27.86	ng	-0.01
23) Nitrobenzene-d5	6.92	82	63355	16.37	ng	0.00
41) 2,4,6-Tribromophenol	10.16	330	15713	21.91	ng	0.00
44) 2-Fluorobiphenyl	8.72	172	108141	17.21	ng	0.00
78) Terphenyl-d14	12.42	244	68194	16.20	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.12	163	29563	4.73	ng	# 97
51) Acenaphthene	9.40	154	11296	2.09	ng	91
70) Phenanthrene	10.86	178	224282	28.20	ng	97
71) Anthracene	10.91	178	53116	6.50	ng	100
74) Fluoranthene	12.03	202	214515	24.91	ng	89
77) Pyrene	12.25	202	189467m	26.10	ng	
80) Benzo(a)anthracene	13.44	228	78461	12.57	ng	96
82) Chrysene	13.47	228	61960m	11.07	ng	
85) Indeno(1,2,3-cd)pyrene	15.82	276	30103	7.33	ng	# 80
87) Benzo(b)fluoranthene	14.43	252	65338m	9.81	ng	
88) Benzo(k)fluoranthene	14.43	252	31886m	5.77	ng	
89) Benzo(a)pyrene	14.71	252	55474	9.83	ng	# 83
91) Benzo(g,h,i)perylene	16.13	276	25360	6.09	ng	# 74

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

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