

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
 Data File : BF081040.D
 Acq On : 18 Aug 2015 10:34
 Operator : UM/IZ
 Sample : G3289-05
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RAMPB-T(0-2)-3

Manual Integrations
 APPROVED

MMDadoda
 8/18/2015 5:14:18 PM

Quant Time: Aug 18 12:57:04 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 17:57:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	61178	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	226281	20.00	ng	0.01
38) Acenaphthene-d10	9.66	164	96777	20.00	ng	0.00
63) Phenanthrene-d10	11.13	188	172700	20.00	ng	0.00
75) Chrysene-d12	13.76	240	132413	20.00	ng	0.01
86) Perylene-d12	15.11	264	85579	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	225720	55.49	ng	0.01
7) Phenol-d6	6.27	99	348399	65.65	ng	0.01
23) Nitrobenzene-d5	7.20	82	249455	50.36	ng	0.01
41) 2,4,6-Tribromophenol	10.44	330	63342	75.51	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	417146	56.48	ng	0.00
78) Terphenyl-d14	12.72	244	324004	52.46	ng	0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.39	163	120183	15.08	ng	100
70) Phenanthrene	11.15	178	45973	4.49	ng	99
74) Fluoranthene	12.33	202	66942	6.06	ng	91
77) Pyrene	12.56	202	61183	5.68	ng	97
80) Benzo(a)anthracene	13.75	228	29702	3.34	ng	98
82) Chrysene	13.78	228	27358m	3.42	ng	
87) Benzo(b)fluoranthene	14.74	252	28821m	4.76	ng	
89) Benzo(a)pyrene	15.05	252	17949	3.40	ng	# 91
91) Benzo(g,h,i)perylene	16.71	276	10364	2.49	ng	# 76

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

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