

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090215\
 Data File : BF081375.D
 Acq On : 3 Sep 2015 4:59
 Operator : UM/IZ
 Sample : G3491-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-EF-01-004-ABCD

Manual Integrations
 APPROVED

MMDadoda
 9/3/2015 1:45:59 PM

Quant Time: Sep 03 05:28:52 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 17:35:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.39	152	65401	20.00	ng	0.00
21) Naphthalene-d8	7.68	136	244869	20.00	ng	0.00
38) Acenaphthene-d10	9.42	164	111294	20.00	ng	0.00
63) Phenanthrene-d10	10.88	188	188685	20.00	ng	0.00
75) Chrysene-d12	13.49	240	106160	20.00	ng	-0.01
86) Perylene-d12	14.80	264	87873	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.96	112	653865	155.12	ng	0.03
7) Phenol-d6	6.07	99	780072	139.81	ng	0.01
23) Nitrobenzene-d5	6.97	82	579594	114.20	ng	0.00
41) 2,4,6-Tribromophenol	10.20	330	141832	143.13	ng	0.00
44) 2-Fluorobiphenyl	8.76	172	845454	97.35	ng	0.00
78) Terphenyl-d14	12.47	244	484103	106.15	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.16	163	126145	14.59	ng	Qvalue # 96
70) Phenanthrene	10.90	178	41813	3.99	ng	97
74) Fluoranthene	12.08	202	62066	5.47	ng	87
77) Pyrene	12.30	202	52355	6.66	ng	97
80) Benzo(a)anthracene	13.47	228	23688	3.50	ng	93
82) Chrysene	13.51	228	17567m	2.90	ng	
85) Indeno(1,2,3-cd)pyrene	15.89	276	10363	2.33	ng	# 93
87) Benzo(b)fluoranthene	14.47	252	22615m	3.60	ng	
88) Benzo(k)fluoranthene	14.48	252	10428m	2.00	ng	
89) Benzo(a)pyrene	14.75	252	16731	3.15	ng	# 88
91) Benzo(g,h,i)perylene	16.22	276	9548	2.44	ng	# 79

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

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