

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100515\
 Data File : BF082061.D
 Acq On : 5 Oct 2015 23:30
 Operator : UM/IZ
 Sample : G3891-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-8(0-2)

Manual Integrations
 APPROVED

MMDadoda
 10/6/2015 2:55:00 PM

Quant Time: Oct 06 06:46:23 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 05 17:58:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	92504	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	367504	20.00	ng	0.00
38) Acenaphthene-d10	9.93	164	153941	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	248605	20.00	ng	0.00
75) Chrysene-d12	14.07	240	196795	20.00	ng	0.01
86) Perylene-d12	15.53	264	134161	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	686589	134.74	ng	0.01
7) Phenol-d6	6.53	99	972292	151.63	ng	0.01
23) Nitrobenzene-d5	7.45	82	530904	96.30	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	183823	117.91	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	1008810	118.60	ng	0.00
78) Terphenyl-d14	13.01	244	737955	154.49	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.65	163	231392	21.78	ng	# 97
70) Phenanthrene	11.44	178	52941	2.87	ng	# 94
74) Fluoranthene	12.63	202	127846	9.60	ng	90
77) Pyrene	12.86	202	117768m	10.02	ng	
80) Benzo(a)anthracene	14.06	228	75139	7.09	ng	# 85
82) Chrysene	14.09	228	90313m	7.54	ng	
85) Indeno(1,2,3-cd)pyrene	16.97	276	41902	5.05	ng	# 77
87) Benzo(b)fluoranthene	15.11	252	99083m	12.94	ng	
88) Benzo(k)fluoranthene	15.12	252	33488m	4.77	ng	
89) Benzo(a)pyrene	15.46	252	58575	8.82	ng	# 87
91) Benzo(g,h,i)perylene	17.41	276	45345	7.94	ng	# 78

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

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