

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021715\
 Data File : BE089603.D
 Acq On : 17 Feb 2015 21:02
 Operator : TP/IZ
 Sample : G1258-01RE
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-1RE

Quant Time: Feb 18 14:39:37 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE021715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 18 14:11:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.35	152	75655	5.00	ng	0.02
7) Naphthalene-d8	12.25	136	375479	5.00	ng	0.01
13) Acenaphthene-d10	16.44	164	191696	5.00	ng	0.02
17) Phenanthrene-d10	19.80	188	457217	5.00	ng	0.03
20) Chrysene-d12	23.66	240	324608	5.00	ng	0.00
27) Perylene-d12	25.34	264	292299	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.50	112	114336	5.81	ng	0.00
4) Phenol-d6	8.63	99	108351	4.18	ng	0.00
8) Nitrobenzene-d5	10.62	82	195783	9.42	ng	0.01
14) 2,4,6-Tribromophenol	18.35	330	95541	16.90	ng	0.03
15) 2-Fluorobiphenyl	14.90	172	494603	9.65	ng	0.02
22) Terphenyl-d14	22.31	244	620353	14.16	ng	0.00
Target Compounds						Qvalue
23) Benzo(a)anthracene	23.64	228	193609m	0.70	ng	
25) Chrysene	23.69	228	175734m	0.55	ng	
26) Indeno(1,2,3-cd)pyrene	26.75	276	3880	0.26	ng	# 96
28) Benzo(b)fluoranthene	24.92	252	17377	0.50	ng	# 91
29) Benzo(k)fluoranthene	24.94	252	9981m	0.18	ng	
30) Benzo(a)pyrene	25.27	252	10649	0.40	ng	# 89
31) Dibenzo(a,h)anthracene	26.77	278	2287	0.27	ng	# 71

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

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