

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE022515\
 Data File : BE089688.D
 Acq On : 25 Feb 2015 20:11
 Operator : TP/JJ
 Sample : G1353-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-2

Quant Time: Feb 26 06:00:57 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE022415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 26 00:03:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	43711	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	204174	5.00	ng	0.00
8) Acenaphthene-d10	16.41	164	94212	5.00	ng	0.00
13) Phenanthrene-d10	19.76	188	161123	5.00	ng	0.00
19) Chrysene-d12	23.65	240	119018	5.00	ng	0.00
25) Perylene-d12	25.34	264	98623	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) Nitrobenzene-d5	10.60	82	75945	6.32	ng	0.00
9) 2-Fluorobiphenyl	14.87	172	171477	8.18	ng	0.00
21) Terphenyl-d14	22.30	244	150920	10.81	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
6) Naphthalene	12.29	128	10485	0.27	ng	# 96
7) 2-Methylnaphthalene	13.94	142	29876	1.33	ng	97
10) Acenaphthylene	16.06	152	41338	0.31	ng	# 76
11) Acenaphthene	16.49	154	8092	0.38	ng	# 44
12) Fluorene	17.74	166	16902	0.71	ng	# 81
14) Hexachlorobenzene	18.99	284	1078	0.26	ng	# 40
15) Pentachlorophenol	19.43	266	693	1.58	ng	93
16) Phenanthrene	19.81	178	44399	1.29	ng	99
17) Anthracene	19.90	178	11747m	0.40	ng	
18) Fluoranthene	21.72	202	9707	0.34	ng	99
20) Pyrene	22.04	202	11815m	0.41	ng	
22) Benzo(a)anthracene	23.64	228	25442	0.33	ng	97
23) Chrysene	23.68	228	33071	0.31	ng	97
24) Indeno(1,2,3-cd)pyrene	26.73	276	13915	0.52	ng	# 54
26) Benzo(b)fluoranthene	24.91	252	2466	0.47	ng	# 92
27) Benzo(k)fluoranthene	24.94	252	7284m	0.36	ng	
28) Benzo(a)pyrene	25.27	252	2825	0.43	ng	# 89
29) Dibenzo(a,h)anthracene	26.76	278	2212	0.48	ng	93
30) Benzo(g,h,i)perylene	27.15	276	15919	0.43	ng	96

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

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