

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020315\
 Data File : BE089477.D
 Acq On : 3 Feb 2015 23:29
 Operator : TP/IZ
 Sample : G1110-08MSD
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-48MSD

Quant Time: Feb 04 02:42:07 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-BE020315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 04 00:53:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.53	152	185585	5.00	ng	0.00
7) Naphthalene-d8	12.44	136	742593	5.00	ng	0.00
13) Acenaphthene-d10	16.62	164	347532	5.00	ng	0.01
17) Phenanthrene-d10	19.94	188	586623	5.00	ng	0.00
20) Chrysene-d12	23.80	240	530871	5.00	ng	0.02
27) Perylene-d12	25.50	264	496297	5.00	ng	0.00
System Monitoring Compounds						
3) 2-Fluorophenol	6.73	112	228993	3.82	ng	0.00
4) Phenol-d6	8.85	99	186177	2.47	ng	0.00
8) Nitrobenzene-d5	10.80	82	460593	8.20	ng	0.02
14) 2,4,6-Tribromophenol	18.53	330	147046	9.77	ng	0.02
15) 2-Fluorobiphenyl	15.07	172	671028	7.33	ng	0.00
22) Terphenyl-d14	22.43	244	720706	9.58	ng	0.00
Target Compounds						Qvalue
2) n-Nitrosodimethylamine	3.53	42	104283	2.82	ng	# 77
5) Phenol	8.88	94	199179	2.42	ng	# 65
6) bis(2-Chloroethyl)ether	9.01	93	417774	7.01	ng	# 98
9) Nitrobenzene	10.85	77	499223	7.77	ng	99
10) 2,4-Dimethylphenol	11.79	122	274825	5.91	ng	97
11) 2,4-Dichlorophenol	12.17	162	244477	7.12	ng	86
12) Hexachlorobutadiene	12.81	225	150960	6.58	ng	99
16) 2,4-Dinitrophenol	16.86	184	126050	14.58	ng	# 79
18) Hexachlorobenzene	19.18	284	237657	7.83	ng	# 82
19) Pentachlorophenol	19.62	266	231535	15.65	ng	97
21) Benzidine	22.12	184	17	0.09	ng	# 1
23) Benzo(a)anthracene	23.78	228	3961738	8.40	ng	98
24) 3,3'-Dichlorobenzidine	23.77	252	27097	0.72	ng	# 92
25) Chrysene	23.83	228	3653614	7.43	ng	94
26) Indeno(1,2,3-cd)pyrene	26.99	276	1141538	8.22	ng	99
28) Benzo(b)fluoranthene	25.06	252	949018	8.35	ng	97
29) Benzo(k)fluoranthene	25.09	252	974109	7.57	ng	99
30) Benzo(a)pyrene	25.43	252	917406	8.27	ng	97
31) Dibenzo(a,h)anthracene	27.02	278	1006169	8.55	ng	96

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

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