

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090115\
 Data File : BE090678.D
 Acq On : 1 Sep 2015 22:21
 Operator : UM/IZ
 Sample : G3462-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MW-2

Manual Integrations
 APPROVED

MMDadoda
 9/2/2015 4:07:59 PM

Quant Time: Sep 02 11:48:40 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 11:40:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	46704	5.00	ng	0.00
6) Naphthalene-d8	10.32	136	200446	5.00	ng	0.00
12) Acenaphthene-d10	14.17	164	105470	5.00	ng	0.00
18) Phenanthrene-d10	16.91	188	277511	5.00	ng	0.00
25) Chrysene-d12	21.07	240	351139	5.00	ng	0.00
32) Perylene-d12	23.36	264	318665	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	16794	1.86	ng	0.00
5) Phenol-d6	6.76	99	13050	1.00	ng	0.00
7) Nitrobenzene-d5	8.72	82	34084	2.57	ng	0.01
13) 2,4,6-Tribromophenol	15.68	330	20740	6.72	ng	0.00
14) 2-Fluorobiphenyl	12.80	172	82780	2.83	ng	0.00
27) Terphenyl-d14	19.53	244	216453	5.57	ng	0.00
Target Compounds						Qvalue
16) Acenaphthene	14.23	154	1052	0.04	ng	90
17) Fluorene	15.24	166	1358	0.04	ng	# 91
22) Phenanthrene	16.95	178	25947	0.38	ng	99
23) Anthracene	17.05	178	6225	0.11	ng	96
24) Fluoranthene	18.96	202	65390	0.90	ng	99
26) Pyrene	19.32	202	61188	0.84	ng	96
28) Benzo(a)anthracene	21.06	228	30182	0.38	ng	98
29) Chrysene	21.11	228	30895	0.31	ng	98
30) Bis(2-ethylhexyl)phthalate	21.01	149	4579	0.29	ng	95
31) Indeno(1,2,3-cd)pyrene	25.73	276	13597	0.16	ng	92
33) Benzo(b)fluoranthene	22.67	252	32730	0.47	ng	98
34) Benzo(k)fluoranthene	22.71	252	15500m	0.19	ng	
35) Benzo(a)pyrene	23.26	252	19614	0.32	ng	95
36) Dibenzo(a,h)anthracene	25.73	278	3019	0.18	ng	# 83
37) Benzo(g,h,i)perylene	26.45	276	15163	0.21	ng	95

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

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