

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
 Data File : BE089966.D
 Acq On : 27 Jun 2015 8:45
 Operator : TP/IZ
 Sample : G2782-09
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X3SS0220006-150624

Manual Integrations
 APPROVED

apatel
 6/29/2015 5:03:29 PM

Quant Time: Jun 29 01:17:22 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 16:51:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	27831	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	135585	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	82161	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	189539	5.00	ng	0.00
25) Chrysene-d12	21.13	240	203701	5.00	ng	0.00
32) Perylene-d12	23.47	264	184318	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	74863	14.90	ng	0.00
5) Phenol-d6	6.81	99	121136	15.73	ng	0.00
7) Nitrobenzene-d5	8.77	82	81989	8.12	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	21397	14.50	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	189628	7.92	ng	0.00
27) Terphenyl-d14	19.60	244	270201	7.68	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.44	128	764	0.03	ng	# 79
11) 2-Methylnaphthalene	12.05	142	625	0.03	ng	# 92
15) Acenaphthylene	13.95	152	5170	0.04	ng	99
16) Acenaphthene	14.30	154	812	0.04	ng	98
17) Fluorene	15.29	166	1197	0.04	ng	# 87
20) Hexachlorobenzene	16.30	284	1192	0.04	ng	# 87
21) Pentachlorophenol	16.63	266	1104	1.00	ng	# 90
22) Phenanthrene	17.01	178	3788m	0.08	ng	
23) Anthracene	17.12	178	1928	0.04	ng	99
24) Fluoranthene	19.02	202	6880	0.13	ng	99
26) Pyrene	19.38	202	6974	0.14	ng	# 96
28) Benzo(a)anthracene	21.12	228	5970	0.12	ng	97
29) Chrysene	21.17	228	6231	0.12	ng	96
30) Bis(2-ethylhexyl)phthalate	21.07	149	5694	1.37	ng	97
31) Indeno(1,2,3-cd)pyrene	25.88	276	3736	0.10	ng	# 100
33) Benzo(b)fluoranthene	22.76	252	6242	0.16	ng	# 90
34) Benzo(k)fluoranthene	22.80	252	3872	0.09	ng	# 75
35) Benzo(a)pyrene	23.36	252	4461	0.13	ng	# 90
36) Dibenzo(a,h)anthracene	25.90	278	1831	0.06	ng	# 71
37) Benzo(g,h,i)perylene	26.62	276	3592	0.10	ng	# 88

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

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