

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
 Data File : BE089964.D
 Acq On : 27 Jun 2015 7:33
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Jun 29 01:12:37 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	23842	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	109197	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	68234	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	165190	5.00	ng	0.00
25) Chrysene-d12	21.13	240	175831	5.00	ng	0.00
32) Perylene-d12	23.47	264	149003	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	3425	0.80	ng	0.00
5) Phenol-d6	6.81	99	5992	0.91	ng	0.00
7) Nitrobenzene-d5	8.77	82	6611	0.81	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	674	0.85	ng	0.00
14) 2-Fluorobiphenyl	12.87	172	17860	0.90	ng	0.00
27) Terphenyl-d14	19.60	244	27131	0.89	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.26	88	2975	1.03	ng	94
3) n-Nitrosodimethylamine	3.55	42	3642	0.85	ng	92
8) Nitrobenzene	8.81	77	7086	0.82	ng	99
9) Naphthalene	10.44	128	22617	0.93	ng	100
10) Hexachlorobutadiene	10.74	225	3571	0.93	ng	99
11) 2-Methylnaphthalene	12.05	142	14982	0.90	ng	99
15) Acenaphthylene	13.95	152	109242	0.91	ng	100
16) Acenaphthene	14.30	154	16624	0.95	ng	98
17) Fluorene	15.29	166	21789	0.99	ng	97
19) 4-Bromophenyl-phenylether	16.18	248	5776	0.86	ng	98
20) Hexachlorobenzene	16.30	284	27089	0.92	ng	99
21) Pentachlorophenol	16.63	266	3501	2.48	ng	91
22) Phenanthrene	17.01	178	37803	0.91	ng	99
23) Anthracene	17.12	178	33412	0.87	ng	99
24) Fluoranthene	19.02	202	39618	0.88	ng	100
26) Pyrene	19.38	202	41751	0.94	ng	100
28) Benzo(a)anthracene	21.12	228	37840	0.90	ng	98
29) Chrysene	21.17	228	40016	0.87	ng	98
30) Bis(2-ethylhexyl)phthalate	21.07	149	3310	0.92	ng	99
31) Indeno(1,2,3-cd)pyrene	25.89	276	31416	0.93	ng	# 100
33) Benzo(b)fluoranthene	22.75	252	30175	0.96	ng	100
34) Benzo(k)fluoranthene	22.80	252	34138	0.96	ng	99
35) Benzo(a)pyrene	23.36	252	25539	0.91	ng	99
36) Dibenzo(a,h)anthracene	25.90	278	25076	0.96	ng	99
37) Benzo(g,h,i)perylene	26.62	276	26134	0.93	ng	99

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

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