

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
 Data File : BE089987.D
 Acq On : 29 Jun 2015 17:21
 Operator : TP/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.5

Quant Time: Jun 30 06:06:59 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 06:02:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	24782	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	104415	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	62646	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	153525	5.00	ng	0.00
25) Chrysene-d12	21.13	240	189974	5.00	ng	0.00
32) Perylene-d12	23.46	264	185618	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	2609	0.56	ng	0.00
5) Phenol-d6	6.80	99	3623	0.51	ng	0.00
7) Nitrobenzene-d5	8.76	82	3805	0.48	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	761	0.80	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	8934	0.48	ng	0.00
27) Terphenyl-d14	19.59	244	15222	0.46	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.25	88	1704	0.50	ng	95
3) n-Nitrosodimethylamine	3.53	42	2256	0.50	ng	# 94
8) Nitrobenzene	8.80	77	4046	0.48	ng	99
9) Naphthalene	10.43	128	11273	0.48	ng	99
10) Hexachlorobutadiene	10.72	225	1780	0.48	ng	99
11) 2-Methylnaphthalene	12.04	142	7432	0.46	ng	98
15) Acenaphthylene	13.95	152	52473	0.47	ng	100
16) Acenaphthene	14.30	154	7851	0.48	ng	99
17) Fluorene	15.29	166	10597	0.48	ng	100
19) 4-Bromophenyl-phenylether	16.18	248	3177	0.50	ng	97
20) Hexachlorobenzene	16.30	284	13206	0.48	ng	100
21) Pentachlorophenol	16.63	266	1088	0.86	ng	99
22) Phenanthrene	17.01	178	19072	0.49	ng	100
23) Anthracene	17.10	178	17113	0.47	ng	100
24) Fluoranthene	19.02	202	20784	0.49	ng	100
26) Pyrene	19.38	202	21450	0.45	ng	99
28) Benzo(a)anthracene	21.11	228	22284	0.48	ng	100
29) Chrysene	21.16	228	23576	0.47	ng	97
30) Bis(2-ethylhexyl)phthalate	21.07	149	5423	0.96	ng	100
31) Indeno(1,2,3-cd)pyrene	25.87	276	21821	0.57	ng	100
33) Benzo(b)fluoranthene	22.75	252	18635	0.47	ng	99
34) Benzo(k)fluoranthene	22.79	252	21667	0.48	ng	99
35) Benzo(a)pyrene	23.35	252	17204	0.48	ng	98
36) Dibenzo(a,h)anthracene	25.89	278	17883	0.53	ng	97
37) Benzo(g,h,i)perylene	26.60	276	18309	0.51	ng	99

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

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