

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082515\
 Data File : BE090627.D
 Acq On : 25 Aug 2015 23:23
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Aug 26 03:28:52 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 14 16:52:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	21831	5.00	ng	0.00
6) Naphthalene-d8	10.33	136	100049	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	57786	5.00	ng	0.00
18) Phenanthrene-d10	16.93	188	147638	5.00	ng	-0.02
25) Chrysene-d12	21.09	240	168450	5.00	ng	0.00
32) Perylene-d12	23.39	264	152622	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	4695	0.74	ng	0.00
5) Phenol-d6	6.78	99	6287	0.68	ng	0.00
7) Nitrobenzene-d5	8.73	82	7024	0.88	ng	0.00
13) 2,4,6-Tribromophenol	15.71	330	1909	0.86	ng	0.00
14) 2-Fluorobiphenyl	12.82	172	15982	0.93	ng	0.00
27) Terphenyl-d14	19.56	244	27255	0.88	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.19	88	3074	1.18	ng	97
3) n-Nitrosodimethylamine	3.49	42	3828	1.02	ng	# 79
8) Nitrobenzene	8.77	77	7308	0.88	ng	97
9) Naphthalene	10.38	128	22232	0.98	ng	99
10) Hexachlorobutadiene	10.67	225	3469	1.04	ng	99
11) 2-Methylnaphthalene	12.01	142	13731	0.93	ng	96
15) Acenaphthylene	13.91	152	98943	0.94	ng	100
16) Acenaphthene	14.25	154	15584	0.98	ng	# 81
17) Fluorene	15.26	166	20001	1.03	ng	99
19) 4-Bromophenyl-phenylether	16.14	248	5459	0.90	ng	84
20) Hexachlorobenzene	16.25	284	26011	0.97	ng	96
21) Pentachlorophenol	16.61	266	1855	0.85	ng	96
22) Phenanthrene	16.98	178	36775	0.95	ng	100
23) Anthracene	17.07	178	33501	0.94	ng	99
24) Fluoranthene	18.98	202	40714	0.99	ng	99
26) Pyrene	19.34	202	43053	0.89	ng	99
28) Benzo(a)anthracene	21.08	228	40587	0.92	ng	97
29) Chrysene	21.13	228	46737	1.03	ng	99
30) Bis(2-ethylhexyl)phthalate	21.02	149	14701	0.68	ng	99
31) Indeno(1,2,3-cd)pyrene	25.78	276	39391	0.92	ng	# 92
33) Benzo(b)fluoranthene	22.69	252	33965	0.94	ng	98
34) Benzo(k)fluoranthene	22.74	252	42573	1.04	ng	98
35) Benzo(a)pyrene	23.29	252	31986	0.95	ng	# 97
36) Dibenzo(a,h)anthracene	25.80	278	29714	0.91	ng	96
37) Benzo(g,h,i)perylene	26.51	276	33229	0.93	ng	96

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

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