

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082515\
 Data File : BE090625.D
 Acq On : 25 Aug 2015 22:11
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Aug 26 03:28:34 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	7.57	152	21608	5.00	ng	0.00
6) Naphthalene-d8	10.33	136	98902	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	58546	5.00	ng	0.00
18) Phenanthrene-d10	16.93	188	150743	5.00	ng	-0.02
25) Chrysene-d12	21.09	240	172480	5.00	ng	0.00
32) Perylene-d12	23.39	264	153687	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	4746	0.76	ng	0.00
5) Phenol-d6	6.78	99	6358	0.69	ng	0.00
7) Nitrobenzene-d5	8.72	82	7364	0.93	ng	0.00
13) 2,4,6-Tribromophenol	15.71	330	1980	0.88	ng	0.00
14) 2-Fluorobiphenyl	12.82	172	16301	0.93	ng	0.00
27) Terphenyl-d14	19.55	244	28526	0.90	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.19	88	2990	1.16	ng	98
3) n-Nitrosodimethylamine	3.49	42	3856	1.03	ng	# 80
8) Nitrobenzene	8.77	77	7324	0.90	ng	98
9) Naphthalene	10.38	128	21958	0.98	ng	99
10) Hexachlorobutadiene	10.67	225	3431	1.04	ng	99
11) 2-Methylnaphthalene	12.01	142	13888	0.95	ng	99
15) Acenaphthylene	13.91	152	100494	0.94	ng	100
16) Acenaphthene	14.26	154	15913	0.99	ng	# 81
17) Fluorene	15.24	166	20742	1.05	ng	99
19) 4-Bromophenyl-phenylether	16.15	248	5514	0.90	ng	78
20) Hexachlorobenzene	16.25	284	27175	1.00	ng	97
21) Pentachlorophenol	16.61	266	1949	0.86	ng	91
22) Phenanthrene	16.98	178	37543	0.95	ng	100
23) Anthracene	17.07	178	34604	0.95	ng	100
24) Fluoranthene	18.98	202	42003	1.00	ng	99
26) Pyrene	19.34	202	44782	0.91	ng	100
28) Benzo(a)anthracene	21.07	228	41563	0.92	ng	100
29) Chrysene	21.13	228	47357	1.02	ng	99
30) Bis(2-ethylhexyl)phthalate	21.02	149	15746	0.70	ng	98
31) Indeno(1,2,3-cd)pyrene	25.79	276	39122	0.90	ng	# 92
33) Benzo(b)fluoranthene	22.69	252	33945	0.93	ng	98
34) Benzo(k)fluoranthene	22.74	252	40974	0.99	ng	98
35) Benzo(a)pyrene	23.29	252	31993	0.94	ng	# 96
36) Dibenzo(a,h)anthracene	25.80	278	29499	0.90	ng	96
37) Benzo(g,h,i)perylene	26.51	276	32941	0.92	ng	96

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

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