

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082515\
 Data File : BE090633.D
 Acq On : 26 Aug 2015 3:01
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Aug 26 04:31:39 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	21337	5.00	ng	0.00
6) Naphthalene-d8	10.33	136	96847	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	56743	5.00	ng	0.00
18) Phenanthrene-d10	16.94	188	147993	5.00	ng	0.00
25) Chrysene-d12	21.09	240	168667	5.00	ng	0.00
32) Perylene-d12	23.39	264	151301	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	4561	0.74	ng	0.00
5) Phenol-d6	6.78	99	6320	0.70	ng	0.00
7) Nitrobenzene-d5	8.72	82	7052	0.91	ng	0.00
13) 2,4,6-Tribromophenol	15.71	330	1813	0.83	ng	0.00
14) 2-Fluorobiphenyl	12.82	172	15481	0.91	ng	0.00
27) Terphenyl-d14	19.56	244	27054	0.88	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.19	88	3056	1.20	ng	95
3) n-Nitrosodimethylamine	3.49	42	3792	1.03	ng	# 81
8) Nitrobenzene	8.77	77	7086	0.89	ng	98
9) Naphthalene	10.38	128	21378	0.97	ng	98
10) Hexachlorobutadiene	10.67	225	3373	1.04	ng	100
11) 2-Methylnaphthalene	12.01	142	13331	0.93	ng	97
15) Acenaphthylene	13.91	152	96506	0.93	ng	100
16) Acenaphthene	14.25	154	15272	0.98	ng	82
17) Fluorene	15.26	166	20166	1.06	ng	99
19) 4-Bromophenyl-phenylether	16.14	248	5518	0.91	ng	85
20) Hexachlorobenzene	16.25	284	26169	0.98	ng	96
21) Pentachlorophenol	16.61	266	1871	0.85	ng	90
22) Phenanthrene	16.98	178	36409	0.94	ng	100
23) Anthracene	17.07	178	33575	0.94	ng	100
24) Fluoranthene	18.98	202	40453	0.98	ng	99
26) Pyrene	19.34	202	42846	0.89	ng	100
28) Benzo(a)anthracene	21.08	228	41621	0.94	ng	97
29) Chrysene	21.13	228	45220	1.00	ng	99
30) Bis(2-ethylhexyl)phthalate	21.02	149	14059	0.66	ng	99
31) Indeno(1,2,3-cd)pyrene	25.78	276	38666	0.91	ng	# 92
33) Benzo(b)fluoranthene	22.70	252	34769	0.97	ng	99
34) Benzo(k)fluoranthene	22.74	252	40410	1.00	ng	98
35) Benzo(a)pyrene	23.29	252	31192	0.93	ng	97
36) Dibenzo(a,h)anthracene	25.80	278	28959	0.90	ng	97
37) Benzo(g,h,i)perylene	26.52	276	32303	0.91	ng	96

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

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