

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080515\
 Data File : BE090364.D
 Acq On : 5 Aug 2015 18:05
 Operator : TP/UM
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.8

Manual Integrations
 APPROVED

MMDadoda
 8/10/2015 10:10:23 AM

Quant Time: Aug 06 02:06:11 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 01:43:02 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	39958	5.00	ng	0.00
6) Naphthalene-d8	10.34	136	177686	5.00	ng	0.00
12) Acenaphthene-d10	14.20	164	102329	5.00	ng	0.00
18) Phenanthrene-d10	16.95	188	239769	5.00	ng	0.00
25) Chrysene-d12	21.10	240	295454	5.00	ng	0.00
32) Perylene-d12	23.40	264	275966	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	4846	0.66	ng	0.00
5) Phenol-d6	6.77	99	8413	0.89	ng	0.00
7) Nitrobenzene-d5	8.73	82	9979	0.94	ng	0.00
13) 2,4,6-Tribromophenol	15.69	330	1390	1.09	ng	-0.02
14) 2-Fluorobiphenyl	12.82	172	23774	0.81	ng	0.00
27) Terphenyl-d14	19.56	244	36719	0.69	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.20	88	4582	0.82	ng	99
3) n-Nitrosodimethylamine	3.49	42	5186	0.78	ng	# 91
8) Nitrobenzene	8.77	77	10526	0.96	ng	100
9) Naphthalene	10.39	128	32917	0.86	ng	98
10) Hexachlorobutadiene	10.69	225	5720	1.08	ng	99
11) 2-Methylnaphthalene	12.01	142	19052	0.81	ng	99
15) Acenaphthylene	13.91	152	138141	0.78	ng	100
16) Acenaphthene	14.26	154	21340	0.83	ng	97
17) Fluorene	15.26	166	28333	0.89	ng	98
19) 4-Bromophenyl-phenylether	16.14	248	7909	0.88	ng	99
20) Hexachlorobenzene	16.27	284	40240	1.13	ng	99
21) Pentachlorophenol	16.61	266	1312	0.91	ng	96
22) Phenanthrene	16.98	178	51074	0.92	ng	100
23) Anthracene	17.07	178	42075	0.86	ng	99
24) Fluoranthene	18.99	202	50228	0.91	ng	100
26) Pyrene	19.35	202	53161	0.69	ng	100
28) Benzo(a)anthracene	21.08	228	55566	0.97	ng	99
29) Chrysene	21.13	228	64849	0.92	ng	99
30) Bis(2-ethylhexyl)phthalate	21.03	149	11643	0.76	ng	99
31) Indeno(1,2,3-cd)pyrene	25.79	276	45558	1.35	ng	97
33) Benzo(b)fluoranthene	22.70	252	40555	1.02	ng	98
34) Benzo(k)fluoranthene	22.75	252	64615m	0.87	ng	
35) Benzo(a)pyrene	23.30	252	37424	0.99	ng	100
36) Dibenzo(a,h)anthracene	25.80	278	34001	1.33	ng	98
37) Benzo(g,h,i)perylene	26.52	276	41250	0.98	ng	99

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

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