

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080515\
 Data File : BE090367.D
 Acq On : 5 Aug 2015 19:53
 Operator : TP/UM
 Sample : SSTDICC005
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

MMDadoda
 8/10/2015 10:10:36 AM

Quant Time: Aug 06 02:12:53 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	7.58	152	37908	5.00	ng	0.00
6) Naphthalene-d8	10.34	136	176832	5.00	ng	0.00
12) Acenaphthene-d10	14.20	164	104620	5.00	ng	0.00
18) Phenanthrene-d10	16.94	188	244698	5.00	ng	0.00
25) Chrysene-d12	21.10	240	294174	5.00	ng	0.00
32) Perylene-d12	23.40	264	268841	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	31118	4.46	ng	0.00
5) Phenol-d6	6.77	99	55058	6.14	ng	-0.01
7) Nitrobenzene-d5	8.72	82	67304	6.36	ng	-0.01
13) 2,4,6-Tribromophenol	15.69	330	13402	6.70	ng	-0.02
14) 2-Fluorobiphenyl	12.82	172	143257	4.75	ng	0.00
27) Terphenyl-d14	19.56	244	229271	4.31	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.19	88	21953	4.99	ng	98
3) n-Nitrosodimethylamine	3.49	42	30452	4.81	ng	# 89
8) Nitrobenzene	8.76	77	73084	6.71	ng	95
9) Naphthalene	10.39	128	178123	4.67	ng	98
10) Hexachlorobutadiene	10.69	225	28929	5.47	ng	100
11) 2-Methylnaphthalene	12.01	142	119450	5.12	ng	98
15) Acenaphthylene	13.91	152	873482	4.83	ng	100
16) Acenaphthene	14.26	154	123659	4.68	ng	90
17) Fluorene	15.26	166	163538	5.01	ng	99
19) 4-Bromophenyl-phenylether	16.14	248	48547	5.29	ng	99
20) Hexachlorobenzene	16.27	284	220344	6.04	ng	98
21) Pentachlorophenol	16.60	266	14849	5.77	ng	90
22) Phenanthrene	16.98	178	279548	4.94	ng	99
23) Anthracene	17.07	178	264420	5.28	ng	100
24) Fluoranthene	18.98	202	314135	5.55	ng	99
26) Pyrene	19.34	202	335210	4.37	ng	99
28) Benzo(a)anthracene	21.08	228	330782	5.80	ng	99
29) Chrysene	21.13	228	367280	5.52	ng	99
30) Bis(2-ethylhexyl)phthalate	21.03	149	121330	4.74	ng	99
31) Indeno(1,2,3-cd)pyrene	25.79	276	330535	7.27	ng	96
33) Benzo(b)fluoranthene	22.70	252	298481	5.67	ng	98
34) Benzo(k)fluoranthene	22.75	252	377835m	5.22	ng	
35) Benzo(a)pyrene	23.30	252	273048	5.63	ng	99
36) Dibenzo(a,h)anthracene	25.81	278	262379	7.09	ng	97
37) Benzo(g,h,i)perylene	26.52	276	284207	5.65	ng	97

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

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