

Data Path : Z:\HPCHEM1\BNA E\DATA\BE092517\
 Data File : BE094112.D
 Acq On : 25 Sep 2017 21:08
 Operator : SJ/JU
 Sample : I5340-12MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 LT-TS-012MS

Manual Integrations
 APPROVED

Sohil
 9/26/2017 5:03:48 PM

Quant Time: Sep 26 02:50:24 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 25 17:37:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1562	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	7386	0.40	ng	0.00
13) Acenaphthene-d10	14.51	164	4866	0.40	ng	0.00
19) Phenanthrene-d10	17.24	188	18461	0.40	ng	0.00
27) Chrysene-d12	21.40	240	15838	0.40	ng	0.00
34) Perylene-d12	23.91	264	15379	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.50	112	850	0.14	ng	0.00
5) Phenol-d6	7.09	99	1343	0.16	ng	0.00
8) Nitrobenzene-d5	9.06	82	1054	0.15	ng	0.00
11) 2-Methylnaphthalene-d10	12.28	152	1652	0.15	ng	0.00
14) 2,4,6-Tribromophenol	16.00	330	364	0.10	ng	0.00
15) 2-Fluorobiphenyl	13.13	172	2160	0.13	ng	0.00
25) Fluoranthene-d10	19.27	212	22892	0.13	ng	0.00
29) Terphenyl-d14	19.85	244	3633	0.13	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.41	88	587	0.193	ng	# 44
3) n-Nitrosodimethylamine	3.73	42	542	0.139	ng	# 90
6) bis(2-Chloroethyl)ether	7.33	93	1000	0.157	ng	99
9) Naphthalene	10.74	128	13376	0.160	ng	99
10) Hexachlorobutadiene	11.03	225	405	0.130	ng	98
12) 2-Methylnaphthalene	12.35	142	2231	0.165	ng	96
16) Acenaphthylene	14.23	152	5714	0.234	ng	98
17) Acenaphthene	14.57	154	2637	0.173	ng	97
18) Fluorene	15.55	166	3816	0.184	ng	96
20) 4-Bromophenyl-phenylether	16.42	248	1246	0.165	ng	97
21) Hexachlorobenzene	16.55	284	835	0.175	ng	98
22) Pentachlorophenol	16.91	266	894	0.204	ng	# 73
23) Phenanthrene	17.27	178	29872	0.632	ng	98
24) Anthracene	17.37	178	10330	0.208	ng	97
26) Fluoranthene	19.30	202	52170	0.986	ng	99
28) Pyrene	19.66	202	53667	1.069	ng	# 96
30) Benzo(a)anthracene	21.38	228	28100	0.527	ng	96
31) Chrysene	21.44	228	31366	0.609	ng	97
32) Bis(2-ethylhexyl)phthalate	21.29	149	165736	1.082	ng	99
33) Indeno(1,2,3-cd)pyrene	26.55	276	19598	0.356	ng	98
35) Benzo(b)fluoranthene	23.14	252	34818	0.627	ng	# 94
36) Benzo(k)fluoranthene	23.18	252	17155m	0.323	ng	
37) Benzo(a)pyrene	23.79	252	26221	0.523	ng	# 94
38) Dibenzo(a,h)anthracene	26.56	278	8279	0.174	ng	# 83
39) Benzo(g,h,i)perylene	27.37	276	18810	0.400	ng	97

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

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