

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011320\
 Data File : BE101044.D
 Acq On : 13 Jan 2020 18:48
 Operator : JU
 Sample : L1103-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SB-1(4-6)

Manual Integrations
 APPROVED

mohammad
 1/14/2020 12:24:05 PM

Quant Time: Jan 14 09:58:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 08 18:36:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	3130	0.40	ng	0.00
7) Naphthalene-d8	10.51	136	13770	0.40	ng	0.01
13) Acenaphthene-d10	14.36	164	9310	0.40	ng	0.00
19) Phenanthrene-d10	17.09	188	21403	0.40	ng	0.00
27) Chrysene-d12	21.27	240	19695	0.40	ng	-0.01
34) Perylene-d12	23.69	264	21903	0.40	ng	-0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.34	112	1458	0.21	ng	0.00
5) Phenol-d6	6.91	99	1799	0.20	ng	0.00
8) Nitrobenzene-d5	8.89	82	1479	0.15	ng	0.01
11) 2-Methylnaphthalene-d10	12.11	152	5080	0.19	ng	0.00
14) 2,4,6-Tribromophenol	15.85	330	315	0.12	ng	0.00
15) 2-Fluorobiphenyl	13.00	172	8278	0.23	ng	0.00
25) Fluoranthene-d10	19.12	212	61283	0.21	ng	-0.01
29) Terphenyl-d14	19.73	244	11717	0.24	ng	-0.01

Target Compounds						Qvalue
6) bis(2-Chloroethyl)ether	7.14	93	28	0.003	ng	# 1
9) Naphthalene	10.55	128	2106	0.014	ng	# 79
10) Hexachlorobutadiene	10.89	225	6	0.001	ng	# 18
12) 2-Methylnaphthalene	12.18	142	263	0.011	ng	# 61
16) Acenaphthylene	14.08	152	625	0.016	ng	# 95
17) Acenaphthene	14.43	154	207	0.008	ng	# 89
18) Fluorene	15.41	166	424	0.012	ng	# 83
20) 4-Bromophenyl-phenylether	16.33	248	11	0.001	ng	# 64
21) Hexachlorobenzene	16.41	284	15	0.001	ng	# 36
23) Phenanthrene	17.13	178	1873	0.032	ng	# 96
26) Fluoranthene	19.15	202	3644	0.050	ng	# 97
28) Pyrene	19.51	202	3989	0.054	ng	# 99
30) Benzo(a)anthracene	21.26	228	2324	0.041	ng	# 81
31) Chrysene	21.30	228	2884	0.035	ng	# 91
32) Bis(2-ethylhexyl)phthalate	21.20	149	10405	0.111	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.23	276	2499	0.038	ng	# 80
35) Benzo(b)fluoranthene	22.95	252	3094	0.050	ng	# 72
36) Benzo(k)fluoranthene	23.00	252	3347m	0.038	ng	#
37) Benzo(a)pyrene	23.58	252	1675	0.028	ng	# 56
38) Dibenzo(a,h)anthracene	26.26	278	811	0.015	ng	# 23
39) Benzo(g,h,i)perylene	27.01	276	3004m	0.045	ng	#

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

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