

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102618\
 Data File : BN003324.D
 Acq On : 27 Oct 2018 02:35
 Operator : MJ/SJ
 Sample : J5521-02MS 10X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 AOC6-01A

Quant Time: Oct 27 05:49:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN102618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 26 15:17:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	12289	0.40	ng	0.00
7) Naphthalene-d8	10.66	136	51147	0.40	ng	-0.01
13) Acenaphthene-d10	14.49	164	29401	0.40	ng	-0.01
19) Phenanthrene-d10	17.24	188	73882	0.40	ng	0.00
27) Chrysene-d12	21.42	240	84373	0.40	ng	0.00
34) Perylene-d12	23.74	264	81809	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.42	112	895	0.03	ng	0.00
5) Phenol-d6	7.01	99	1147	0.03	ng	0.00
8) Nitrobenzene-d5	9.02	82	732	0.03	ng	0.00
11) 2-Methylnaphthalene-d10	12.25	152	2223	0.03	ng	0.00
14) 2,4,6-Tribromophenol	15.98	330	296	0.03	ng	0.00
15) 2-Fluorobiphenyl	13.12	172	3318	0.03	ng	-0.01
25) Fluoranthene-d10	19.26	212	4667	0.02	ng	0.00
29) Terphenyl-d14	19.86	244	5736	0.03	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.34	88	514	0.032	ng	# 80
3) n-Nitrosodimethylamine	3.67	42	142	0.019	ng	# 1
6) bis(2-Chloroethyl)ether	7.29	93	836	0.024	ng	96
9) Naphthalene	10.71	128	3464	0.028	ng	# 93
10) Hexachlorobutadiene	10.99	225	491	0.021	ng	# 96
12) 2-Methylnaphthalene	12.32	142	2350	0.027	ng	97
16) Acenaphthylene	14.21	152	4348	0.033	ng	98
17) Acenaphthene	14.56	154	2075	0.023	ng	95
18) Fluorene	15.54	166	3032	0.027	ng	100
20) 4-Bromophenyl-phenylether	16.43	248	917	0.021	ng	# 86
21) Hexachlorobenzene	16.54	284	964	0.021	ng	99
22) Pentachlorophenol	16.88	266	245	0.020	ng	# 85
23) Phenanthrene	17.28	178	7793	0.038	ng	98
24) Anthracene	17.37	178	4442	0.026	ng	99
26) Fluoranthene	19.29	202	13926	0.060	ng	99
28) Pyrene	19.66	202	13769	0.059	ng	98
30) Benzo(a)anthracene	21.40	228	10563	0.048	ng	96
31) Chrysene	21.45	228	11298	0.047	ng	97
32) Bis(2-ethylhexyl)phthalate	21.32	149	3557	0.054	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.09	276	9828	0.040	ng	96
35) Benzo(b)fluoranthene	23.03	252	14403	0.057	ng	# 87
36) Benzo(k)fluoranthene	23.08	252	8277	0.033	ng	# 75
37) Benzo(a)pyrene	23.63	252	10594	0.046	ng	# 86
38) Dibenzo(a,h)anthracene	26.11	278	5663	0.026	ng	# 78
39) Benzo(g,h,i)perylene	26.82	276	9541	0.044	ng	93

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

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