

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN022718\
 Data File : BN000181.D
 Acq On : 27 Feb 2018 09:16
 Operator : JU/SJ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

Sohil
 2/28/2018 5:12:07 PM

Quant Time: Feb 27 14:56:06 2018
 Quant Method : Z:\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 14:08:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	6042	0.40	ng	0.00
7) Naphthalene-d8	11.29	136	22293	0.40	ng	0.00
13) Acenaphthene-d10	15.06	164	11048	0.40	ng	0.00
19) Phenanthrene-d10	17.78	188	25445	0.40	ng	0.00
27) Chrysene-d12	21.90	240	20732	0.40	ng	0.00
34) Perylene-d12	24.37	264	16932	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.93	112	1299	0.10	ng	0.00
5) Phenol-d6	7.58	99	1584	0.10	ng	0.00
8) Nitrobenzene-d5	9.63	82	1293	0.10	ng	0.00
11) 2-Methylnaphthalene-d10	12.85	152	3308	0.10	ng	0.00
14) 2,4,6-Tribromophenol	16.53	330	308	0.08	ng	0.00
15) 2-Fluorobiphenyl	13.69	172	5227	0.10	ng	0.00
25) Fluoranthene-d10	19.77	212	6489	0.10	ng	0.00
29) Terphenyl-d14	20.33	244	3732	0.10	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.66	88	608	0.086	ng	94
3) n-Nitrosodimethylamine	4.05	42	368m	0.079	ng	
6) bis(2-Chloroethyl)ether	7.85	93	2029	0.099	ng	97
9) Naphthalene	11.34	128	6519	0.100	ng	95
10) Hexachlorobutadiene	11.62	225	1115	0.101	ng	97
12) 2-Methylnaphthalene	12.92	142	4113	0.096	ng	99
16) Acenaphthylene	14.78	152	4694	0.090	ng	100
17) Acenaphthene	15.13	154	3932	0.096	ng	95
18) Fluorene	16.09	166	4902	0.096	ng	# 80
20) 4-Bromophenyl-phenylether	16.97	248	1351	0.092	ng	# 81
21) Hexachlorobenzene	17.09	284	1946	0.099	ng	95
22) Pentachlorophenol	17.43	266	633	0.121	ng	92
23) Phenanthrene	17.82	178	8712	0.097	ng	96
24) Anthracene	17.91	178	5972	0.085	ng	96
26) Fluoranthene	19.80	202	8145	0.088	ng	# 96
28) Pyrene	20.15	202	8595	0.106	ng	98
30) Benzo(a)anthracene	21.87	228	6347	0.097	ng	98
31) Chrysene	21.93	228	7653	0.100	ng	99
32) Bis(2-ethylhexyl)phthalate	21.76	149	2967	0.122	ng	# 97
33) Indeno(1,2,3-cd)pyrene	26.91	276	6104	0.086	ng	# 90
35) Benzo(b)fluoranthene	23.61	252	7538	0.098	ng	94
36) Benzo(k)fluoranthene	23.66	252	6783	0.090	ng	# 72
37) Benzo(a)pyrene	24.26	252	5709	0.091	ng	# 69
38) Dibenzo(a,h)anthracene	26.93	278	4354	0.078	ng	# 78
39) Benzo(g,h,i)perylene	27.70	276	6108	0.092	ng	# 81

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

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