

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091823\
 Data File : BN027732.D
 Acq On : 18 Sep 2023 13:34
 Operator : MA/JU
 Sample : SSTDICC3.2
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC3.2

Quant Time: Sep 18 14:07:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 18 14:03:35 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	1793	0.400	ng	0.00
7) Naphthalene-d8	10.831	136	5735	0.400	ng	0.00
13) Acenaphthene-d10	14.641	164	3978	0.400	ng	-0.01
19) Phenanthrene-d10	17.393	188	10079	0.400	ng	# 0.00
29) Chrysene-d12	21.578	240	12114	0.400	ng	# 0.00
35) Perylene-d12	24.075	264	12454	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.546	112	13041	2.764	ng	0.00
5) Phenol-d6	7.156	99	17132	3.003	ng	0.00
8) Nitrobenzene-d5	9.209	82	14655	3.461	ng	0.00
11) 2-Methylnaphthalene-d10	12.411	152	27655	3.271	ng	0.00
14) 2,4,6-Tribromophenol	16.140	330	5508	3.584	ng	-0.01
15) 2-Fluorobiphenyl	13.272	172	44398	2.983	ng	0.00
27) Fluoranthene-d10	19.421	212	90312	3.517	ng	0.00
31) Terphenyl-d14	20.006	244	68614	2.821	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.422	88	6287	2.872	ng	98
3) n-Nitrosodimethylamine	3.769	42	7784	3.250	ng	# 98
6) bis(2-Chloroethyl)ether	7.445	93	17336	3.108	ng	99
9) Naphthalene	10.885	128	49822	3.150	ng	95
10) Hexachlorobutadiene	11.109	225	8327	2.886	ng	# 99
12) 2-Methylnaphthalene	12.487	142	36383	3.272	ng	97
16) Acenaphthylene	14.373	152	56571	3.138	ng	99
17) Acenaphthene	14.705	154	37428	3.085	ng	99
18) Fluorene	15.693	166	54671	3.326	ng	99
20) 4,6-Dinitro-2-methylph...	16.847	198	512	3.471	ng	# 27
21) 4-Bromophenyl-phenylether	16.574	248	17194	3.230	ng	# 80
22) Hexachlorobenzene	16.661	284	18712	2.985	ng	98
23) Atrazine	16.847	200	14060	3.484	ng	# 93
24) Pentachlorophenol	17.033	266	9435	3.796	ng	98
25) Phenanthrene	17.430	178	95711	3.215	ng	100
26) Anthracene	17.530	178	86636	3.407	ng	100
28) Fluoranthene	19.449	202	125074	3.514	ng	99
30) Pyrene	19.816	202	129629	2.727	ng	99
32) Benzo(a)anthracene	21.560	228	138564	3.285	ng	99
33) Chrysene	21.614	228	142747	3.104	ng	100
34) Bis(2-ethylhexyl)phtha...	21.435	149	64121	3.217	ng	98
36) Indeno(1,2,3-cd)pyrene	26.694	276	196092	3.770	ng	100
37) Benzo(b)fluoranthene	23.297	252	152133	3.148	ng	# 91
38) Benzo(k)fluoranthene	23.347	252	157826	3.175	ng	# 91
39) Benzo(a)pyrene	23.958	252	150751	3.346	ng	# 87
40) Dibenzo(a,h)anthracene	26.718	278	160893	3.919	ng	94
41) Benzo(g,h,i)perylene	27.519	276	166336	3.533	ng	98

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

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