

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110724\
 Data File : BN034885.D
 Acq On : 07 Nov 2024 10:02
 Operator : RC/JU
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.1

Quant Time: Nov 07 14:39:35 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN110724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 14:34:20 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	5582	0.400	ng	0.00
7) Naphthalene-d8	10.340	136	16157	0.400	ng	0.00
13) Acenaphthene-d10	14.208	164	7091	0.400	ng	0.00
19) Phenanthrene-d10	16.952	188	14302	0.400	ng	# 0.00
29) Chrysene-d12	21.149	240	7364	0.400	ng	0.00
35) Perylene-d12	23.318	264	6365	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.192	112	1587	0.102	ng	0.00
5) Phenol-d6	6.752	99	2007	0.097	ng	0.00
8) Nitrobenzene-d5	8.707	82	1268	0.101	ng	0.00
11) 2-Methylnaphthalene-d10	11.935	152	2127	0.097	ng	0.00
14) 2,4,6-Tribromophenol	15.698	330	148	0.101	ng	0.00
15) 2-Fluorobiphenyl	12.829	172	3107	0.104	ng	0.00
27) Fluoranthene-d10	18.990	212	2865	0.089	ng	0.00
31) Terphenyl-d14	19.603	244	1450	0.105	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.184	88	786	0.111	ng	96
3) n-Nitrosodimethylamine	3.487	42	973	0.102	ng	96
6) bis(2-Chloroethyl)ether	7.012	93	1836	0.103	ng	99
9) Naphthalene	10.383	128	4482	0.100	ng	97
10) Hexachlorobutadiene	10.682	225	738	0.103	ng	# 100
12) 2-Methylnaphthalene	12.011	142	2614	0.095	ng	98
16) Acenaphthylene	13.919	152	3250	0.095	ng	99
17) Acenaphthene	14.272	154	2291	0.097	ng	98
18) Fluorene	15.255	166	2861	0.097	ng	99
21) 4-Bromophenyl-phenylether	16.158	248	759	0.100	ng	# 84
22) Hexachlorobenzene	16.269	284	936	0.102	ng	98
23) Atrazine	16.431	200	457	0.083	ng	# 85
25) Phenanthrene	16.989	178	4356	0.099	ng	99
26) Anthracene	17.088	178	3451	0.091	ng	99
28) Fluoranthene	19.018	202	3977	0.086	ng	100
30) Pyrene	19.385	202	4013	0.108	ng	99
32) Benzo(a)anthracene	21.131	228	2670	0.093	ng	97
33) Chrysene	21.185	228	3005	0.099	ng	96
34) Bis(2-ethylhexyl)phtha...	21.095	149	1811	0.110	ng	# 99
36) Indeno(1,2,3-cd)pyrene	25.473	276	2880	0.102	ng	98
37) Benzo(b)fluoranthene	22.672	252	2681	0.096	ng	# 62
38) Benzo(k)fluoranthene	22.713	252	2806	0.096	ng	# 56
39) Benzo(a)pyrene	23.222	252	2082	0.094	ng	# 50
40) Dibenzo(a,h)anthracene	25.491	278	2210	0.101	ng	# 84
41) Benzo(g,h,i)perylene	26.128	276	2359	0.101	ng	91

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

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