

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060121\
 Data File : BN014817.D
 Acq On : 01 Jun 2021 17:30
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN060121

Manual Integrations
 APPROVED

mohammad
 6/2/2021 10:53:17 AM

Quant Time: Jun 01 18:09:13 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN060121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 01 17:40:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.717	152	618	0.40	ng	0.00
7) Naphthalene-d8	10.492	136	2144	0.40	ng	0.00
13) Acenaphthene-d10	14.359	164	1499	0.40	ng	0.01
19) Phenanthrene-d10	17.104	188	3687	0.40	ng	# 0.00
29) Chrysene-d12	21.311	240	4893	0.40	ng	# 0.00
35) Perylene-d12	23.576	264	4944	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.317	112	697	0.43	ng	0.00
5) Phenol-d6	6.895	99	849	0.40	ng	0.00
8) Nitrobenzene-d5	8.857	82	674m	0.37	ng	0.00
11) 2-Methylnaphthalene-d10	12.093	152	1506	0.37	ng	0.00
14) 2,4,6-Tribromophenol	15.856	330	285	0.36	ng	0.00
15) 2-Fluorobiphenyl	12.976	172	2187	0.40	ng	0.00
27) Fluoranthene-d10	19.151	212	4806	0.35	ng	0.00
31) Terphenyl-d14	19.752	244	4379	0.38	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.271	88	352	0.34	ng	# 88
3) n-Nitrosodimethylamine	3.658	42	52	0.29	ng	# 1
6) bis(2-Chloroethyl)ether	7.145	93	678	0.39	ng	95
9) Naphthalene	10.543	128	2319	0.39	ng	96
10) Hexachlorobutadiene	10.835	225	648	0.40	ng	# 94
12) 2-Methylnaphthalene	12.169	142	1583	0.38	ng	98
16) Acenaphthylene	14.067	152	2100	0.37	ng	97
17) Acenaphthene	14.422	154	1644	0.39	ng	98
18) Fluorene	15.412	166	2196	0.38	ng	99
20) 4,6-Dinitro-2-methylph...	15.501	198	123	0.41	ng	# 81
21) 4-Bromophenyl-phenylether	16.313	248	896	0.37	ng	90
22) Hexachlorobenzene	16.422	284	1097	0.38	ng	98
23) Atrazine	16.580	200	491	0.32	ng	99
24) Pentachlorophenol	16.763	266	174	0.42	ng	# 87
25) Phenanthrene	17.152	178	3922	0.38	ng	99
26) Anthracene	17.238	178	2984	0.35	ng	97
28) Fluoranthene	19.181	202	5118	0.36	ng	99
30) Pyrene	19.543	202	5181	0.37	ng	99
32) Benzo(a)anthracene	21.300	228	4989	0.36	ng	99
33) Chrysene	21.343	228	6364	0.38	ng	97
34) Bis(2-ethylhexyl)phtha...	21.226	149	1471	0.38	ng	99
36) Indeno(1,2,3-cd)pyrene	25.852	276	7485	0.37	ng	98
37) Benzo(b)fluoranthene	22.894	252	6841m	0.38	ng	
38) Benzo(k)fluoranthene	22.939	252	7084	0.37	ng	98
39) Benzo(a)pyrene	23.476	252	6101	0.39	ng	# 90
40) Dibenzo(a,h)anthracene	25.872	278	6132	0.36	ng	99
41) Benzo(g,h,i)perylene	26.550	276	6860	0.37	ng	100

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

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