

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040824\
 Data File : BN030781.D
 Acq On : 08 Apr 2024 22:39
 Operator : MA/JU
 Sample : PB160073BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB160073BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 04/10/2024
 Supervised By :mohammad ahmed 04/10/2024

Quant Time: Apr 09 04:27:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN033024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 00:04:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.682	152	6208	0.400	ng	-0.02
7) Naphthalene-d8	10.464	136	19279	0.400	ng	-0.02
13) Acenaphthene-d10	14.327	164	12177	0.400	ng	-0.01
19) Phenanthrene-d10	17.078	188	30448	0.400	ng	-0.01
29) Chrysene-d12	21.267	240	26720	0.400	ng	-0.02
35) Perylene-d12	23.504	264	22488	0.400	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.285	112	5926	0.438	ng	-0.01
5) Phenol-d6	6.859	99	6542	0.376	ng	-0.01
8) Nitrobenzene-d5	8.830	82	4083	0.301	ng	-0.02
11) 2-Methylnaphthalene-d10	12.058	152	11050m	0.375	ng	-0.02
14) 2,4,6-Tribromophenol	15.824	330	1831	0.398	ng	-0.01
15) 2-Fluorobiphenyl	12.948	172	13984	0.297	ng	-0.02
27) Fluoranthene-d10	19.108	212	25503	0.311	ng	-0.02
31) Terphenyl-d14	19.711	244	21252	0.344	ng	-0.02
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.219	88	2285	0.305	ng	# 29
3) n-Nitrosodimethylamine	3.530	42	1940	0.268	ng	# 88
6) bis(2-Chloroethyl)ether	7.119	93	5478	0.319	ng	96
9) Naphthalene	10.517	128	16438	0.308	ng	100
10) Hexachlorobutadiene	10.795	225	3133	0.287	ng	# 99
12) 2-Methylnaphthalene	12.134	142	10935	0.312	ng	100
16) Acenaphthylene	14.038	152	16057	0.291	ng	99
17) Acenaphthene	14.381	154	10991	0.288	ng	99
18) Fluorene	15.375	166	15375	0.309	ng	100
20) 4,6-Dinitro-2-methylph...	15.460	198	820	0.347	ng	# 90
21) 4-Bromophenyl-phenylether	16.271	248	4966	0.270	ng	# 79
22) Hexachlorobenzene	16.370	284	6124	0.279	ng	99
23) Atrazine	16.556	200	3289	0.246	ng	99
24) Pentachlorophenol	16.718	266	1763	0.366	ng	98
25) Phenanthrene	17.115	178	27239	0.302	ng	97
26) Anthracene	17.202	178	21470	0.283	ng	99
28) Fluoranthene	19.135	202	31661	0.295	ng	100
30) Pyrene	19.502	202	34834	0.333	ng	100
32) Benzo(a)anthracene	21.249	228	28196	0.298	ng	99
33) Chrysene	21.303	228	29613	0.292	ng	99
34) Bis(2-ethylhexyl)phtha...	21.187	149	14995	0.365	ng	# 99
36) Indeno(1,2,3-cd)pyrene	25.764	276	25068	0.252	ng	99
37) Benzo(b)fluoranthene	22.832	252	30216m	0.330	ng	
38) Benzo(k)fluoranthene	22.872	252	27782	0.303	ng	97
39) Benzo(a)pyrene	23.407	252	23118	0.316	ng	# 90
40) Dibenzo(a,h)anthracene	25.782	278	19414	0.246	ng	97
41) Benzo(g,h,i)perylene	26.454	276	21330	0.251	ng	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040824\
 Data File : BN030781.D
 Acq On : 08 Apr 2024 22:39
 Operator : MA/JU
 Sample : PB160073BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :

BNA_N

ClientSampleId :

PB160073BS

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 04/10/2024

Supervised By :mohammad ahmed 04/10/2024

Quant Time: Apr 09 04:27:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN033024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 00:04:36 2024
 Response via : Initial Calibration

